Introduction to the "Limited Site Investigation Report Seattle Public Utilities North Recycling and Disposal Station" conducted by AMEC Earth & Environmental, July 8, 2008

What's the Background on the Site?

- The City of Seattle has operated a 4.27-acre solid waste transfer station in the Fremont/Wallingford area since 1967, collecting recyclables and solid waste dropped off by Seattle households, businesses, and collection trucks. A neighboring property, the Oroweat Baking Company site at 1550 N. 34th Street, was bought by the City in 2005 to support modernizing the transfer station facilities.
- For many years, vehicles have entered the site and unloaded their solid waste into a concrete enclosed area below grade. Recyclables are stored at the ground surface and shipped off to recycling facilities. Waste materials are removed from the collection area daily and moved offsite for disposal. Annually the City removes the asphalt layer over the concrete floor of the collection area and disposes of it, laying a new and clean asphalt floor.

What are the Plans for the Site?

 Plans are in place to demolish the existing structures and build a new transfer station and recycling facility. The new transfer station would be built to current standards, enclosing operations more completely to minimize odor, dust, and noise and offering opportunities to improve site lighting, landscaping, and traffic management. The majority of recycling, employee, and administrative facilities would be located at the Oroweat site.

What Does the Report Say?

- In this report, the City contracted with an independent environmental firm to sample and analyze soil and groundwater on the NRDS and Oroweat properties to help characterize the degree of any potential contamination that may currently exist from years of solid waste operations or previous site uses. The soil and groundwater samples were analyzed for petroleum hydrocarbons (oils), volatile organic compounds (VOCs), semi volatile organic compounds (SVOCs) metals, pesticides, herbicides, polychlorinated biphenyls, and (PCBs). There were no detections for pesticides, herbicides, or PCBs. The findings regarding petroleum hydrocarbons, VOCs, SVOCs and metals revealed nothing unusual for a site of this type. Highlights of the report include:
 - The study found traces of diesel and oil in soil, with one soil sample at the uphill background location along the north site boundary at a level exceeding state cleanup standards for lubrication oil.
 - Out of about 200 parameters analyzed in groundwater only two parameters (volatile organic compounds [solvents]) were detected

- in one location that exceeded the most stringent state cleanup standards.
- There were no detections of semi volatile organic compounds or metals that exceeded the most stringent state cleanup standards.
- The Oroweat property has long been known to have gasoline and diesel fuel contamination in soil beneath the existing building resulting from truck fueling and storage, which remained after earlier regulatory cleanups of contamination. In addition, during a prior underground storage tank removal and cleanup project, some petroleum contaminated soil was left near a retaining wall by the southern perimeter of the NRDS facility because the soil could not be removed without compromising the stability of the wall.

As indicated in the SEPA checklist and DNS, a plan for removal, treatment, and/or other management of contaminated soil and groundwater on these sites will be developed prior to construction, with appropriate worker safety, management of contaminated material, and equipment decontamination. SPU will work with permitting and regulatory agencies as appropriate.

How Will You Find Out What Happens?

 SPU will keep the community posted through website updates, and emails to our e-mail list.

Talk to Us at Any Time

- If you have questions you can contact Jeff Neuner at (206) 684-7693
- You can also email us at <u>Newstations@Seattle.gov</u>. If you would like to receive any email updates, please send us your email address.
- The webpage with facility updates can be found at www.seattle.gov/util
 Type in "Facility Update" in the search field to find a link to the Facility Update page.



LIMITED SITE INVESTIGATION REPORT

Seattle Public Utilities North Recycling and Disposal Station
1350 and 15050 NNorth 34th Street
Seattle, Washington 98103

Submitted to:

Seattle Public Utilities 707 South Plummer Street Seattle, Washington 98134

Submitted by:

AMEC Earth & Environmental, Inc. 11810 North Creek Parkway North Bothell, Washington 98011 11335 N.E. 122nd Way, Suite 100 Kirkland, Washington 98034

July 8, 2008 revised March 26, 2009

AMEC Project No. 8-915-16341-0



July 8, 2008 revised March 26, 2009 8-915-16341-0

Seattle Public Utilities 707 South Plummer Street Seattle, Washington 98134

Attention: Juan Carlos Ramirez

Subject: Limited Site Investigation Report

Seattle Public Utilities North Recycling and Disposal Station

1350 and 15<u>5</u>00 NNorth 34th Street

Seattle, Washington 98103

Dear Mr. Ramirez:

AMEC Earth & Environmental, Inc. is pleased to submit the results of the Limited Site Investigation study for the above-referenced property located in Seattle, Washington. This report has been prepared for the exclusive use of the Seattle Public Utilities and its attorneys, for specific application to this project, in accordance with generally accepted environmental practices.

We appreciate the opportunity to work with SPU on this project. If you have any questions or desire further information, please feel free to contact the undersigned at (425) 368-1000.

Sincerely,

AMEC Earth & Environmental, Inc.

Cherilyn Inouye Project Manager Meg Strong, LG Senior Associate

Attachments

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EXECUTIVE SUMMARY

AMEC Earth & Environmental, Inc. (AMEC) is pleased to present this <u>revised</u> Limited Site Investigation Report which was requested by the Seattle Public Utilities (SPU) Materials Laboratory. This report replaces the July 8, 2008 version, and revisions are noted in strike-out and underlined text. This report summarizes the results of an environmental investigation conducted at the properties located at 1350 and <u>4500–1550 NNorth</u> 34th Street, Seattle, Washington 98103 (Site). The purpose of this Limited Site Investigation is to conduct a preliminary screening of the Site soil and groundwater to evaluate if these media have been impacted by current and historic activities onsite and evaluate if these contaminants pose a health risk to future construction workers. At the request of the SPU Materials Laboratory, AMEC collected soil and groundwater samples which were submitted for environmental analyses during geotechnical drilling operations performed by SPU during March 2008. This report presents AMEC's observations, subsurface investigation data, laboratory analytical results, and conclusions from this March 2008 investigation.

AMEC understands that SPU is considering expansion options for the North Recycling and Disposal Station (NRDS). Therefore, future potential receptors on the Site would include construction workers and facility operations personnel. The exposure pathways of the Site's future chemicals of concern, and of its current receptors, are: direct contact, inhalation, and ingestion.

The Site is located in the North Lake Union area of Seattle. <u>Land use in t</u>The area surrounding the Site is characterized <u>with as a mix of light industrial</u>, commercial, and residential development. The Site consists of two tax parcels separated by a north-south public road (Carr Place <u>NNorth</u>). Four one-story buildings are present in the western portion of the Site which has been occupied by the SPU NRDS (1350 <u>NNorth</u> 34th Street) since 1967.

A multi-story building is present in the eastern portion of the Site (15950 NNorth 34th Street) which is currently vacant. Oroweat Baking Company formerly owned and occupied this portion of the Site until 1998. Previous environmental investigations conducted on the former Oroweat property indicate that the soil and groundwater on this property have been impacted by a release of petroleum fuel and solvents associated with the fueling and maintenance of trucks on this property.

During this investigation, five soil borings were advanced using a hollow stem auger drill rig. Three of the soil borings were located on the NRDS property, and two of the soil borings were advanced on the former Oroweat property. Soil samples were collected from each of the soil borings with the exception of No soil samples were collected from boring B-2, located on the former Oroweat property. The three soil borings located on the NRDS property were completed as monitoring wells. Groundwater was not encountered during the advancement of the soil borings on the former Oroweat property so no monitoring wells were installed.

Soil borings advanced dDuring this investigation encountered fill materials was observed in the soil borings at from the ground surface ranging in thickness from 0 toto a maximum depth of 7

feet. The fill materials generally consisted of silty sand with gravel. The fill materials are underlain by glacial till which consists of dense to very dense fine to medium sand with silt and gravel.

Groundwater was encountered during drilling at depths between approximately between–9 to 15 feet below ground surface (bgs). Groundwater levels in the three newly installed monitoring wells was approximately 7 feet to the top of casing (TOC) in two wells at the floor level of the transfer station (MW-1 and MW-2) and approximately 13.5 feet to the TOC in MW-3 which is located at the northern edge of the NRDS and approximately 15 feet higher topographically. Groundwater flow is expected to follow the slope of the topography to the southwest.

To evaluate the <u>potential</u> impacts of current and former Site operations to the Site soil and groundwater, at least one soil and groundwater sample was collected from four of the five soil borings advanced on the Site and analyzed for Site chemicals of potential concern (COPCs), which include: petroleum-range hydrocarbons, volatile organic compounds, semi-volatile organic compounds, metals, polychlorinated biphenyls, and pesticides/herbicides. With the exception of a faint petroleum odor observed in soil boring B-3, no soil staining or petroleum odors were observed during this investigation.

Results from this limited Site Investigation indicate that COPCs are present in the Site soil and groundwater. The limited number of COPCs detected in both the soil and groundwater onsite suggests that the effects of each COPC to human health and the environment may be evaluated against the Washington State Department of Ecology (Ecology) Model Toxics Control Act (MTCA) Method A cleanup levels (CULs) for Unrestricted Land Use. For COPCs that have CUL values for both the MTCA Methods A and B, the values listed under Method A were used in this evaluation. If a COPC did not have a Method A CUL value, then the Method B value was used in this evaluation.

In the Site soil, oil-range hydrocarbons was the only COPC to be detected above the MTCA Method A CULs. No COPCs were detected above Method B CULs where there were no Method A values for soil. The elevated concentration of oil range hydrocarbons was reported in one soil sample collected in boring B-5 near the northern property perimeter located approximately 15 feet higher in elevation than the floor level of the main building. Because the sample was collected from a location hydraulically upgradient from the facility's operations, the source of the petroleum may potentially be from an offsite source.

In the Site groundwater, the following chlorinated solvents were detected: tetrachloroethene (PCE), 1,1-dichloroethene (DCE), and 1,2-dichloroethane (DCA). In addition, low levels of gasoline constituents were detected in the central area of the Site. No COPCs were detected above MTCA Method A CULs or Federal/State MCLs. One COPC (1,1-DCE), which did not have a MTCA Method A CUL, was detected in a concentration above the MTCA Method B CUL. No other COPCs were detected above Method B CULs if there were no value for the MTCA Method A CUL.

The PCE was detected in the upgradient monitoring well MW-3 located along the Site's northern perimeter, and the eventual daughter (breakdown) products of PCE (DCE and DCA) were

detected in a monitoring well MW-2 located in the southeastern corner of the 1350 North 34th Street property. The presence of these COPCs in the groundwater indicates that there may have been a release of solvents near or upgradient of the well locations. Low levels of gasoline constituents were also detected in the groundwater in the central area of the Site. These detections indicate that there may have been a localized release of gasoline onsite.

Results from this limited Site Investigation indicate that COPCs are present in the Site groundwater in concentrations below Washington State Department of Ecology (Ecology) Model Toxics Control Act Method A cleanup levels and Federal Maximum Contaminant Levels for drinking water, and oil range hydrocarbons are present in the Site soil in concentrations above Ecology Method A cleanup levels. The elevated concentration of oil range hydrocarbons are present in an upgradient location adjacent to a public street and utility corridor which suggests that the source of the petroleum release may be from an offsite source. In addition, the presence of the COPCs (petroleum compounds, constituents of fuel, and solvents) in the groundwater indicates that there may be a release of these compounds near or upgradient of the soil and groundwater sample locations.

In addition, the 15500 NNorth 34th Street property has a restrictive covenant due to the presence of gasoline and diesel-impacted soil and perched groundwater remaining below the building footprint. If future development plans include demolition of the building or disturbance of the petroleum-impacted soil, Ecology is required to be notified, and the diesel-impacted soil will be required to be managed as a potentially hazardous waste petroleum contaminated soil.

On the basis of the results of this investigation, it is AMEC's opinion that the following actions be conducted onsite:

- 1. If the soil below the 15500 NNorth 34th Street building will be disturbed, Ecology should be notified in accordance with the property's restrictive covenant, and a soil management plan should be prepared. The purposes of the plan is to address would be 1) to address worker safety monitoring and hazard prevention in the known contaminated areas onsite; and 2) to prepare site-specific proposed plans to segregate, manage, and dispose of contaminated soil; and equipment decontamination.
- 2. If the NRDS is to be demolished, it is recommended that additional soil and groundwater samples be collected below building foundations in areas where floor drains are located. Gasoline constituents were detected in the Site groundwater in concentrations below MTCA CULs and State MCLs. The presence of the gasoline constituents in only one of the three monitoring wells indicates that a localized source may be present within the main building. If the NRDS is to be demolished, it is recommended that additional soil and groundwater samples be collected below building foundations in areas where floor drains are located. The presence of the gasoline constituents in the groundwater in the central area indicates that there is a release of gasoline onsite.
- 3. Should development plans include the demolition of the NRDS building located in the southeast corner of the 1350 North 34th Street property, the soil below the building and the electrical subsurface utility line connecting this building to the main building should

be evaluated for the presence of petroleum, specifically gasoline and diesel. A lens of petroleum-impacted soil was reported in these areas in 1994 during the removal of two underground storage tanks. Locate and review all documents pertaining to the former underground storage tanks (USTs)Where practical, the petroleum contaminated soil was removed. However, to maintain the structural integrity of two retaining walls and a building structure, the cleanup activities did not remove all of the petroleum contaminated soil onsite. on the NRDS property and evaluate the extent of the reported petroleum release associated with the USTs. Following this evaluation, the UST area may require additional environmental investigation and remediation to achieve closure from Ecology. This Additional cleanup of the Site soil could be conducted in conjunction with Site redevelopment.

- 4. Collect groundwater <u>samples</u> from the three of the existing wells during the dry season to monitor the presence of VOCs, specifically the solvents and gasoline constituents in the groundwater. The purpose of the groundwater evaluation is to <u>assess the effects of seasonal variation and to confirm</u> that the VOCs, solvents, and gasoline constituents have remained at levels below Ecology <u>MTCA Method A CULscleanup levels</u>.
- 5. Evaluate the extent of the elevated oil-range hydrocarbons detected along the northern property boundary.
- 6. Survey the elevations of the top of casing in the monitoring wells onsite. The groundwater gradient onsite may be calculated following completion of this survey.

The preceding summary is intended for introduction and reference only. A complete reading of this report is recommended.

Seattle Public Utilities Project No.: 8-915-16341-0

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LIMITED SITE INVESTIGATION REPORT NORTH RECYCLING AND DISPOSAL, SEATTLE, WASHINGTON JULY 2008 REVISED MARCH 2009

1.0 INTRODUCTION

AMEC Earth & Environmental, Inc. (AMEC) is pleased to present this <u>revised</u> Limited Site Investigation Report which was requested by the Seattle Public Utilities (SPU) Materials Laboratory. This report replaces the July 8, 2008 version, and revisions are noted in strike-out and underlined text. This report summarizes the results of an environmental investigation conducted at the properties located at 1350 and 15950 North 34th Street, Seattle, Washington 98103 (Site). The purpose of this Limited Site Investigation was to conduct a preliminary screening of the Site soil and groundwater to evaluate if these media have been impacted by current and historic activities onsite and evaluate if these contaminants pose a health risk to future construction workers. At the request of the SPU Materials Laboratory, AMEC collected soil and groundwater samples which were submitted for environmental analyses during geotechnical drilling operations performed by SPU during March 2008. This report presents AMEC's observations, subsurface investigation data, laboratory analytical results, and conclusions from the March 2008 investigation. This report is also a companion document to the Seattle Public Utilities Geotechnical Data Report-Solid Waste Facilities Master Plan (SWFMP): Subsurface Data Collection, Phase 2 North Recycling and Disposal Station (NRDS).

It is AMEC's understanding that the purpose of this initial investigation was to conduct a screening of subsurface Site chemicals of potential concern (COPCs).

Seattle Public Utilities Project No.: 8-915-16341-0

2.0 SITE BACKGROUND

It is the understanding of AMEC that SPU has owned and operated a <u>solid</u> waste transfer station on the 1350 N<u>orth</u> 34th <u>Street</u> property since 1967 and has recently purchased the 15<u>5</u>00 N<u>orth</u> 34th Street property which is located east of the waste transfer station. Both properties have been used for commercial and light industrial purposes since the 19<u>2</u>60's.

2.1 Site Location and Description

The Site is located in the North Lake Union area of Seattle. The Land use in the area surrounding the Site is characterized with as a mix of light industrial, commercial, and residential development.

The Site consists of two adjoining tax parcels (Parcel numbers 2264500450 and 4083306930) and has an approximate area of 5.18 acres. Carr Place North, a north-south public road, separates the two properties. The Site is bordered to the north by North 35th Street followed by residential homes and L&O Distributing in the northwest; to the east by Woodlawn Avenue North followed by residential homes and the Essential Baking Company; to the south by North 34th Street followed by Institute for Systems Biology, Yogi Way, Northwest Lighting Service, JS Jefferson & Plaster, and Impart Media Group; and to the west by the Kite Shop, Subway Restaurant, and Offshore Store (Figure 2).

Four one-story-buildings are present on the 1350 NNorth 34th Street property (North Recycling and Disposal Station [NRDS]). These buildings include:

- a 58,804-square foot <u>buildingwarehouse</u> used to <u>temporarily storetransfer</u> solid waste, (transfer station)_constructed in 1967,
- a 351-square foot equipment shop building constructed in 1967,
- a 600-square foot office constructed in 196794, and
- a 1,200 square foot shed constructed in 1994.

One multi-story building is present on the 15<u>5</u>00 NNorth 34th Street property. It is currently <u>used</u> by SPU for various recycling activities and occasional meetingsvacant and was formerly occupied by Oroweat.

According to topographic data available on the King County Tax Assessor web site, Site elevations range from 75 feet above mean sea level (msl) in the northeast corner of the Site (1500–1550 NNorth 34th Street property) to approximately 30 feet above msl in the western corner of the Site (1350 NNorth 34th Street property) (Figure 3). The topography of the NRDS property (1350 NNorth 34th Street), which most likely originally sloped to the southwest towards Lake Union, has been altered to accommodate construction of the solid waste transfer handling facility. In order to create a flat surface for the building footprint, a portion of the hillside in the eastern half of the property was removed resulting in steep embankments along the northern and eastern edges of the Site. All four sides of this property slope towards the approximate center which represents a topographic depression.

Surface soil is exposed in landscaped areas along the perimeters of both properties and also along the driveway entrances to the NRDS property (1350 NNorth 34th Street). Surface water on the property is anticipated to either infiltrate the unpaved areas or flow overland towards stormwater catch basins onsite and south and southwest of the Site. The stormwater catch basins discharge to the sanitary sewer

2.2 Current Use of Site

SPU has operated a solid waste transfer station on the western portion of the Site, NRDS property (1350 NNorth 34th Street) since 1967. Residential and commercial customers may drop off their solid waste and recyclable materials at the facility. According to Mr. Ken Armstrong the Director of SPU Solid Waste and Field Operations, solid waste and recyclable materials brought to the facility are placed into the main warehousetransfer building "the pit," and large bulky items (e.g., tires and household appliances) are temporarily stored on a paved surface in the northeast portion of the property. The solid waste is removed from the facility daily and disposed of offsite in Oregontransported to a permitted disposal facility, and the recyclableing materials are removed from the facility periodically, dependingent on storage space.

Mr. Armstrong also indicated that a nonprofit agency and a tenant store building materials in the building on the eastern portion of the Site, on the former Oroweat property (15050 North 34th Street). In addition SPU periodically conducts meetings in the office area of the building.

2.3 Previous Environmental Investigations

Prior to the field investigation, AMEC reviewed 7 reports provided by SPU regarding the property located at 4500–1550 NNorth. 34th Street (former Oroweat property—also listed with street address 1550 N 34th—Street). On May 14, 2008, AMEC also reviewed environmental records on file for the Oroweat and NRDS properties at the Washington State Department of Ecology (Ecology) Bellevue office.

A review of the reports and Ecology file indicates that a truck maintenance facility was present on the Oroweat property. Three underground storage tanks (USTs) containing diesel, gasoline and waste oil were installed on the property by 1968. Releases from these USTs were suspected to be the sources of petroleum contamination in soil and groundwater on the property which was first detected in 1992. In addition to petroleum hydrocarbons, chlorinated solvents such as perchloroethylene (PCE) and trichloroethane (TCA), used for parts cleaning during truck maintenance activities, are suspected of being released as these chemicals have also been detected in soil and groundwater on the property.

Several episodes of remediation activities have been conducted on the Oroweat property including the excavation and removal of petroleum-contaminated soils, in-situ contaminated soil treatment by geo-oxidation using electro-chemical oxidation, localized groundwater treatment with hydrogen peroxide and in-situ soil and groundwater treatment with hydrogen peroxide. In March 2001 Ecology issued a No Further Action Letter and Restrictive Covenant for the property. Under the terms of the Restrictive Covenant, Ecology mandated notification in the event that soil or groundwater below the Former Oroweat property was to be disturbed and or

the removal of any contaminated media. In November 2003, Urban Redevelopment LLC submitted a Voluntary Cleanup Program Remedial Closure Report to Ecology followed by quarterly groundwater monitoring reports completed in 2004. The results of the 2004 groundwater monitoring indicated that the <u>detected concentrations of the</u> chemicals evaluated were below Ecology cleanup levels. A letter from Ecology to SPU (February 17, 2005), stated that Ecology did not require additional groundwater monitoring on the Oroweat property. This letter did not modify any other conditions or requirements of Ecology's "no further action" determination, including the Restrictive Covenant filed with King County on March 28, 2001. An updated Restrictive Covenant to reflect the no further action determination for the Site was not in the Ecology file reviewed by AMEC.

AMEC also reviewed Ecology's environmental file for the NRDS property located at 1350 North 34th Street and the *North Recycling and Disposal Station Site Characterization and Cleanup* Report (Herrera 1997) prepared after the decommissioning of two USTs onsite. The Ecology file contained a "UST Closure and Site Assessment Notice" and a "UST Notice of Confirmed Release" filed in 1994. One 10,000-gallon UST formerly containing diesel fuel and one 3,000-gallon UST formerly containing gasoline and later diesel fuel were permanently removed and disposed off-site. Reportedly, the USTs were present onsite in 1960installed in 1967. According to the 1997 cleanup report, During the UST removal, petroleum-impacted soil contamination was observed during the removal of the USTs. The petroleum-impacted soil was excavated to the practical extent possible. However, to maintain the structural integrity of two retaining walls and a building located in the southeast corner of the property, all of the petroleum-impacted soil could not be excavated. Consequently, petroleum-impacted soil was left in-place in the southeast corner of the property. Observations recorded during the cleanup activities described the impacted soil in the southern excavation sidewall as a "2 to 3 inch (gray-stained soil lens) 2 to 3 feet wide."

On June 10, 2002, Ecology requested additional information regarding site cleanup activities from the City of Seattle. The Ecology letter stated that the current status of the Site is "Cleanup Started" and additional information is needed. Ecology was concerned that petroleum hydrocarbon contamination may migrate under the public road (road not specified). No records reporting concentrations of diesel- and oil-range TPHs in soil and groundwater and the extent of contamination were available for AMEC's review at Ecology.

Ecology's files also contained a regular correspondence between Ecology and the City of Seattle pertaining to the solid waste permits and annual reports dated 2005, 2006, and 2007. No violation of solid waste handling at the site was recorded.

2.4 Geology and Hydrogeology

Seattle lies within the Puget Sound Lowland, an elongate structural and topographic basin bordered by the Cascade and Olympic Mountains. The geology of the Seattle area is dominated by a complex, alternating, and incomplete sequence of glacial and interglacial deposits that rest upon an irregular bedrock surface. Post-glacial sediments are poorly consolidated, as much as 300 meters thick in deep alluvial valleys and susceptible to ground failure during earthquakes. (Troost et al, 2003).

Based on recent <u>regional</u> geological mapping, surficial geologic conditions at the Site consist of artificial fill materials overlying Pleistocene glacial drift deposits of the Fraser Glaciation. Glacial drift underlying fill materials at the Site consist of the Vashon till. The artificial fill materials are <u>described asobserved onsite include</u> a mixture of gravel, sand, <u>and silt, concrete, garbage, slag</u> and other materials of substantial areal extent or thickness (>2 meters) placed as a result of <u>human activity</u>. The Vashon Till is described as a compact diamict of silt, sand and subrounded to well-rounded gravel, glacially transported and deposited under ice. Commonly fractured and containing intercalated sand lenses, the Vashon Till generally forms undulating elongated surfaces. The upper 1 meter of the till is generally weathered. The unit is described as dense to medium dense (Troost et al, 2005).

Soil borings advanced during this investigation encountered fill materials at the ground surface ranging in thickness from 0 to 7 feet. The fill materials generally consisted of silty sand with gravel. The fill materials are underlain by glacial till which consists of dense to very dense fine to medium sand with silt and gravel. In addition, an exposure of the Vashon Till was noted in the hillside just north of an electrical transformer in the west central portion of the Site. The outcrop is in an area that was excavated for placement of the transformer.

Groundwater was encountered during drilling at depths between approximately between 9 to 15 feet bgs. Groundwater levels in the three (MW-1, MW-2, and MW-3) newly installed monitoring wells was approximately 7 feet to the top of casing (TOC) in wells at the floor of the transfer station (MW-1 and MW-2) and approximately 13.5 feet to the TOC in MW-3 which is located approximately 15 feet higher topographically. Groundwater flow is expected to follow the slope of the topography to the southwest.

3.0 SOIL AND GROUNDWATER INVESTIGATION

This section discusses the subsurface investigation performed by AMEC between March 10 and 14, 2008. AMEC collected soil and groundwater samples from four of the five soil borings advanced on the Site. Description of the sample collection and field observations are presented below, and the analytical results of the environmental samples are presented in Section 4 of this document.

3.1 Analytical Program for Soil and Groundwater

To evaluate the impact of current and former Site operations to the Site soil and groundwater, at least one soil sample was collected and analyzed from each boring for the Site COPCs which include: petroleum range hydrocarbons, volatile organic compounds (VOCs), semi-VOCs (SVOCs), metals, polychlorinated biphenyls (PCBs), pesticides, and herbicides.

3.2 Soil Sample Collection

Soil samples were collected from 4 soil borings (B-1, B-3, B-4 and B-5) advanced to a maximum depth of 30.3 feet bgs using a hollow stem auger operated by Gregory Drilling, Inc. of Redmond, Washington. A geotechnical boring, B-2, was also advanced on the Site; however environmental samples were not collected from this boring.

As illustrated on Figure 3, the soil borings were spatially distributed evenly across the property and also placed strategically downgradient of current potential source areas. Relatively undisturbed subsurface soil core samples were collected at 2½ and 5-foot intervals by driving a 1½-foot-long, 2-inch-diameter split-spoon sampler. Following sample collection, soil samples were visually inspected and field screened for VOCs using an organic vapor meter (OVM) equipped with a photo-ionization detector (PID). In each soil boring, SPU personnel recorded a description of the soil lithology. AMEC's field observations were added to SPU's soil boring logs which are presented in the SPU Materials Laboratory geotechnical report.

Soil samples submitted to the laboratory were selected on the basis of field observations and soil recovery. One soil sample in each soil boring, except B-2, was selected for laboratory analyses. Due to the large volume of soil required for analyses from each boring, AMEC collected soil samples from multiple intervals. A faint petroleum odor was noted in the 2.5 to 4-foot sample obtained from soil boring, B-3. However, no PID measurements above background levels were observed in any of the borings subjected to the PID soil samples checked by AMEC.

Soil samples were placed into laboratory-prepared containers. Soil samples to be analyzed for volatile petroleum constituents (e.g., total petroleum hydrocarbons [TPH] in the gasoline range and VOCs) were collected using a soil core syringe and inserted into a pre-tared 40-millileter (ml) VOA vial in accordance with EPA Method 5035. Samples to be analyzed for non-volatile analyses (e.g. metals, SVOCs, pesticides/herbicides, and PCBs) were collected in new 4-ounce soil jars. Samples were sealed, labeled, and placed in a cooler with ice for transport and delivery to the analytical laboratory under proper chain-of-custody.

3.3 Monitoring Well Installation and Development

Soil borings B-1, B-4, and B-5 were completed as two-inch diameter monitoring wells (MW-1, MW-2, and MW-3, respectively). Groundwater was not encountered in B-3, therefore a monitoring well was not installed. The monitoring wells were installed to monitor groundwater elevations and to assess potential impacts of contaminated soil on the Site to groundwater beneath the Site.

Ten-foot sections of slotted PVC screen were installed between 10 and 20 feet bgs in MW-1 and MW-3 and between 5 and 15 feet bgs in MW-2. During drilling, groundwater was observed at depths ranging from 9 to 15 feet bgs. A filter pack size compatible with the well screen was installed by gravity in the annulus of each borehole. The filter pack enclosed the screened section of the well and extended at least 1 foot above the top of the well screen. Bentonite chips were placed to form a seal extending from the top of the filter pack to approximately two feet bgs, and a traffic-rated flush mount well cover was placed over the well casing and securely cemented into the ground.

On March 12, 2008 the newly installed wells were developed using a decontaminated surge block and submersible pump to remove fine material and turbid water from the wells. The static water levels in the wells ranged between approximately 5.5 feet and 13 feet bgs-below the TOC during the well development. To develop the well, the surge block was thrust up and down the screened interval for a period of approximately 10 minutes. Following surging, the submersible pump was placed into the well to remove the turbid water. Three cycles of surging and purging were conducted in each well. All three of the wells pumped dry during well development; however, well development continued until three well volumes were removed or the well was purged dry three times. Water quality measurements recorded on field forms during the well development are presented in Appendix A.

3.4 Groundwater Sample Collection

Groundwater samples were collected from the three monitoring wells on March 14, 2008. To ensure that the groundwater samples collected were representative of the aquifer, groundwater from the permanent wells was purged prior to sample collection using a peristaltic pump with disposable tubing. Groundwater was removed at a low flow rate (approximately 300 milliliter per minute [ml/min]). At regular intervals, the purged water was monitored for temperature, pH, specific conductivity, turbidity, dissolved oxygen, and oxidation-reduction potential using a Horiba U-22 water quality meter. Ghese parameters are recorded on groundwater sampling field forms which are included in Appendix B. The well purging was completed upon stabilization of water quality parameters defined as two readings within 10 percent of the previous measurement. It should be noted that the dissolved oxygen meter malfunctioned during the groundwater sample collection. Consequently the dissolved oxygen measurements were not evaluated to determine the stabilization of groundwater parameters.

Groundwater samples to be analyzed for TPH in the gasoline range and VOCs were collected in 40-ml VOA vials pre-preserved with hydrochloric acid; samples to be analyzed for SVOCs, herbicides, pesticides, and PCBs were collected in unpreserved 1-Liter amber bottles; and samples to be analyzed for metals were collected in 500-ml plastic bottles. Immediately

following collection, samples were sealed, labeled, and placed in a cooler with ice for delivery to the analytical laboratory under chain-of-custody protocols. The groundwater samples to be analyzed for metals were filtered in the laboratory and preserved prior to analysis.

3.5 Equipment Decontamination

Following each soil sampling interval, the sampling equipment was cleaned and decontaminated using a scrub brush and non-phosphate detergent solution followed by a deionized-water and isopropyl alcohol rinse followed by air-drying. Decontamination liquids and rinsate were containerized and stored onsite as investigation-derived waste (IDW).

3.6 Investigation Derived Waste

Two waste streams were generated during drilling activities (soil cuttings and decontamination liquids/purge water). The IDW was segregated by matrix, placed into 55-gallon drums, labeled, and stored on the Site in the northeast corner of the NRDS property (1350 NNorth 34th Street). Four soil drums and two liquid waste drums were generated during the March 2008 investigation. On the basis of the analytical results, it is anticipated that this the IDW will be was disposed of as non-hazardous waste.



4.0 SITE INVESTIGATION RESULTS

This section presents the overall findings of the AMEC Site investigation. The results of the laboratory analyses of soil and groundwater samples collected during AMEC's investigation are summarized below and are presented in Tables 1 through 14. All tables are located at the end of this report, and the analytical laboratory data reports are presented in Appendix C. The soil analytical results discussed in this section have been preliminarily screened against values of the Ecology Model Toxics Control Act (MTCA) Method A and B cleanup levels (CULs) for Unrestricted Land Use—and—Method—B—CULs. The groundwater analytical results were preliminarily screened against Federal/State Maximum Contaminant Limits (MCLs) for drinking water,—and Ecology MTCA A and B CULs for Unrestricted Land Use, and MTCA B CULs.

4.1 Soil Analytical Results

The laboratory results for soil samples collected are discussed below and are summarized on Tables 1 to 7. During the March 2008 investigation, soil samples were analyzed for one or more of the following: TPH gasoline by NWTPH-Gx, TPH diesel by NWTPH-Dx, VOCs by EPA Method 8260B, SVOCs by EPA Method 8270D-SIM, metals (arsenic, barium, cadmium, chromium, lead, mercury, selenium, and silver) using EPA 6010/7471, herbicides by EPA Method 8151, pesticides by EPA Method 8081A, and PCBs by EPA Method 8082. The results of these analyses are presented below.

4.1.1 Total Petroleum Hydrocarbons

Four primary and one duplicate soil sample were analyzed for gasoline-, diesel-, and oil-range hydrocarbons using NWTPH Gx and Dx. The results for the TPH <u>analyses</u> are presented on Table 1 <u>and Figure 5</u>. Gasoline-range hydrocarbons were not detected in the soil samples in concentrations above the laboratory detection limits. Diesel range hydrocarbons were detected in only one soil boring, B-1, at a concentration of 150 milligrams per kilogram (mg/kg). Lube oil range hydrocarbons were detected in three primary and one duplicate soil sample. Detections of lube oil ranged from 110 mg/kg in B-4 to 2,700 mg/kg in B-5. Per WAC 173-340, the sum of the diesel and oil-range hydrocarbons exceeded the MTCA Method A CUL of 2,000 mg/kg for diesel and oil in the soil samples collected from B-5 at 7 feet bgs. No other detections of petroleum hydrocarbons exceeded MTCA Method A CULs.

4.1.2 Volatile Organic Compounds

Four primary and one duplicate soil sample were analyzed for VOCs using EPA Method 8260B. The results for the VOCs are presented on Table 2. Three VOCs (acetone, tetrachloroethene PCE, and toluene) were detected in one or more soil samples. None of the VOCs were detected above the MTCA Method A or B CULs.

4.1.3 Semi-Volatile Organic Compounds

Four primary and one duplicate soil sample was analyzed for SVOCs using EPA Method 8270D SIM. The results for the SVOCs are presented on Table 3. Four SVOCs (bis[2-

ethylhexyl]phthalate, di-n-butylphthalate, phenanthrene, and pyrene) were detected in one or more soil samples. None of the detections exceeded the MTCA Method A and B CULs. It should be noted that the laboratory detection limit for n-nitrosodimethylamine (0.035 mg/kg) exceeded the MTCA Method B CUL of 0.02 mg/kg. N-nitrosodimethylamine was not detected in any of the soil samples. Because it is typically associated with rocket fuel, n-nitrosodimethylamine is not anticipated to be present in the soil onsite.

4.1.4 Metals

Four primary and one duplicate soil sample were analyzed for arsenic, barium, cadmium, total chromium, soluble hexavalent chromium, lead, mercury, selenium, and silver using EPA Methods 6010B and 7471A. A summary of the analytical results is presented on Table 4. Barium and total chromium were detected in all of the soil samples. The detected concentrations of these metals were below their respective MTCA Method A and B CULs.

4.1.5 Pesticides

Four soil samples were analyzed for pesticides using EPA Method 8081A, and a summary of the analytical results are presented on Table 5. The pesticide compounds were not detected above laboratory detection limits in these samples.

4.1.6 Herbicides

Four soil samples were analyzed for herbicides using EPA Method 8151, and a summary of the analytical results are presented on Table 6. The herbicide compounds were not detected above laboratory detection limits in these samples.

4.1.7 Polychlorinated Biphenyls

Four soil samples were analyzed for PCBs using EPA Method 8082. A summary of the analytical results are presented on Table 7. The PCB compounds were not detected above laboratory detection limits in these samples.

4.2 Groundwater Analytical Results

The laboratory results for the groundwater samples collected from the three newly installed monitoring wells are discussed below and are summarized on Tables 8 to 14. During the March 2008 investigation, groundwater samples were analyzed for: TPH gasoline by NWTPH-Gx TPH diesel by NWTPH-Dx, VOCs by EPA Method 8260B, SVOCs by EPA Method 8270D-SIM, metals (arsenic, barium, cadmium, chromium, lead, mercury, selenium, and silver) using EPA 6010/7471, herbicides by EPA Method 8151, and pesticides by EPA Method 8081A. The results of these analyses are presented below.

The groundwater analytical results were preliminarily screened against Federal Maximum Contaminant Limits for drinking water, Ecology MTCA A CULs for Unrestricted Land Use, and MTCA B CULs for Unrestricted Land Use.

4.2.1 Total Petroleum Hydrocarbons

Three primary and one duplicate groundwater sample were analyzed for gasoline, diesel, and oil range hydrocarbons using NWTPH Gx and Dx. The results of the TPH analyses are presented on Table 8. Petroleum-range hydrocarbons were not detected in concentrations above the laboratory detection limits in the samples.

4.2.2 Volatile Organic Compounds

Three primary and one duplicate groundwater sample were analyzed for VOCs using EPA Method 8260B. The results of the analyses are presented on Table 9. Ten VOCs were detected in one or more samples. None of the detections exceeded Federal MCLs for drinking water or MTCA Method A CULs (where available). The detections of 1,1-dichloroethene (DCE) in MW-2 (31 µg/L); 1,2-dichloroethane (DCA) in MW-2 (3.4µg/L); and tetrachloroethenePCE in MW-1–3 (0.810 µg/L) exceeded the MTCA Method B CULs for these VOCs (Figure 6). No other detections exceeded the screening criteria.

The laboratory detection limits for three compounds (1,1-dichloroethene DCE, 1,2,3-trichloropropane, and 1,2-dibromo-3-chloropropane) exceeded their respective MTCA Method B CULs. 1,2,3-trichloropropane and 1,2-dibromo-3-chloropropane were not detected above laboratory detection limits in the groundwater samples collected.

4.2.3 Semi-Volatile Organic Compounds

Three primary and one duplicate groundwater sample were analyzed for SVOCs using EPA Method 8270D SIM. The results are presented on Table 10. Two SVOCs (di-n-butylphthalate and pyrene) were detected in one or more groundwater samples. None of the detections exceeded Federal MCLs for drinking water or MTCA Method A and B CULs.

It should be noted that the laboratory method detection limit of the following eight compounds exceeded the MTCA Method B CULs: 1,2-diphenylhydrazine; 3,3'-dichlorobenzidine; benzidine; bis(2-Chloroethyl)ether; hexachlorobenzene; hexachlorobutadiene; n-nitrosodimethylamine; and pentachlorophenol. These compounds were not detected above the laboratory detection limits in any of the groundwater samples.

4.2.4 Metals

Three primary and one duplicate groundwater sample were analyzed for dissolved arsenic, barium, cadmium, chromium, lead, mercury, selenium, and silver using EPA Methods 6010B and 7471A. A summary of the analytical results is presented on Table 11. Barium was detected in the groundwater samples collected from MW-1 and MW-3. The detected concentrations did not exceed Federal MCLs for drinking water or MTCA Methods A and B CULs. No other metal was detected above laboratory detection limits.

4.2.5 Pesticides

Three primary and one duplicate groundwater sample were analyzed for pesticides using EPA Method 8081A, and a summary of the analytical results are presented on Table 12. No pesticide

compounds were detected above the laboratory detection limits. It should be noted that the laboratory detection limit for aldrin exceeded the MTCA Method B CUL.

4.2.6 Herbicides

Three primary and one duplicate groundwater sample were analyzed for herbicides using EPA Method 8151, and a summary of the analytical results are presented on Table 13. No herbicide compounds were detected above the laboratory detection limits.

4.2.7 PCBs

Three primary and one duplicate groundwater sample were analyzed for PCBs using EPA Method 8082. A summary of the analytical results are presented on Table 14. No PCB compounds were detected above the laboratory detection limits.



5.0 QUALITY ANALYSIS / QUALITY CONTROL

A qualitative assessment of the soil and groundwater analytical data from this study is presented in this section. The overall usability of the data was acceptable with respect to the results of the precision, accuracy, representativeness, comparability, and completeness evaluations. Quality control (QC) measures implemented during this study included an assessment of sample handling and management and an evaluation of field and laboratory QC samples.

Field QC samples are collected for laboratory analyses to assess sampling and analytical accuracy and precision, efficiency of sampling equipment decontamination, and potential cross-contamination from the time of sample collection to laboratory analyses. Field QC samples consisted of field duplicates, equipment rinsates, trip blanks, and field blanks. Laboratory QC samples for this investigation consisted of method blanks, laboratory duplicates, laboratory control samples / laboratory control sample duplicate (LCS/LCSD) and matrix spike / matrix spike duplicate (MS/MSD) samples. The 90% project data quality objective for valid measurements was met for this project. In general, it is important to recognize that no analytical data are guaranteed to be correct, even if all QC audits are passed. Strict QC serves to increase confidence in data, but any reported value may potentially contain error.

5.1 Sample Handling and Management

All samples were stored in a cooler with bagged ice and securely maintained in AMEC's possession under chain <u>andof</u> custody protocols. All samples were received at Onsite Environmental in good condition at a cooler temperature of 4°C, which was within the EPA guidelines of ≤6°C. All samples were analyzed within the EPA-recommended maximum holding time of 14 days for soils and preserved samples for VOCs, SVOCs, TPHs, PCBs, pesticides, and herbicides; and within 28 days for metals.

A review of the laboratory data indicates the following data reporting issues at the laboratory:

- Sample B-5_6 had an internal standard for volatile analysis that was outside the control limits. The laboratory states that for B-5_6 the results and the corresponding Practical quantitation limits (PQLs) from bromobenzene onward should be considered estimates. It should be noted that the 19 VOCs following the bromobenzene were not detected above the laboratory detection limits in sample B-5_6, and the PQL values for the 19 compounds are below the MTCA A and B CULs.
- The laboratory reported that sample B-4_6 had interfering compounds that may be masking the presence of acetone.
- The laboratory reported that the toluene in samples B-1_5, B-3_3, B-4-6, B-5_6, and B-5_7 may be present due to cross contamination during storage. In AMEC's professional opinion, the toluene results from samples B-1_5, B-3_3, B-4-6, B-5_6, and B-5_7 should be used with caution due to the absence of other compounds usually associated with toluene contamination.

4,4'-DDT and Methoxychlor recoveries in the continuing calibration verifications (CCV's)
were low; therefore the results in the samples for these compounds are biased low and
maybe greater than reported. These compounds were not detected above the laboratory
detection limits in the soil or groundwater samples.

5.2 Field QC Samples

Field duplicates were collected at a rate of 10% per media (e.g. groundwater and soil). Three duplicate soil samples and two duplicate groundwater samples were collected and analyzed. The field duplicate samples were analyzed using the same analytical program as the primary samples. It should be noted that due to limited sample volume, no duplicate samples were analyzed for herbicides, pesticides, and PCBs. A review of the laboratory data indicates that duplicate pairs showed precision within Quality Analysis Project Plan (QAPP)-specified limits with the following exception.

The field duplicate B-5_7 was collected at the same location as B-5_6. The percent deviation between the acetone and tetrachloroethene PCE recoveries is greater than the acceptance limit of 40% for soils at 63% and 121%. The acetone and tetrachloroethene PCE results should be considered estimated.

One equipment rinsate <u>sample</u>, <u>EB-1</u>, was collected for this field effort. The rinsate was collected by pouring deionized water over a decontaminated split-spoon sampler (soil sampling equipment). The rinsate water was then poured directly into laboratory containers. No target compounds were detected in the soil equipment rinsate sample. Because disposable groundwater sampling (bailers and dedicated tubing) was utilized, no equipment rinsate was collected for the groundwater sampling equipment. <u>No target compounds were detected in the soil equipment rinsate sample blank with the following exception:</u>

• Equipment Blank EB-1 had detections of methylene chloride at 2.5 μg/L, chloroform at 1.7 μg/L, and diethylphthalate 1.9 μg/L. These chemicals were not detected in the associated samples; therefore data usability was not impacted.

Trip blanks and field blanks were collected daily, and following review of the preliminary laboratory data, select trip and field blanks were analyzed for volatile compounds (EPA method 8260B and/or NWTPH Gx). No target compounds were detected in the trip blank samples, or the field blanks with the following exceptions:

- •Equipment Blank, EB-1, had detections of methylene chloride at 2.5 μg/L, chloroform at 1.7 μg/L, and diethylphthalate 1.9 μg/L. The associated samples were non-detected for methylene chloride, chloroform, and diethylphthalate; therefore data usability was not impacted.
- A field blank, Field blank_3, had a detection of chloroform at 1.2 μg/L. Consequently, <u>t</u>The detection of chloroform in sample MW-2 at 0.66 μg/L should be considered non-detect.

• The trip blank, TRIP BLANK 2, was not analyzed for vinyl acetate. This compound was not detected in the associated samples were all non-detect for vinyl acetate; therefore data usability was not impacted.

5.3 Laboratory QC Samples

The samples are processed in batches of less than 20 samples per QC set. The QC sets requirements vary by method, but generally consist of a method blank, laboratory control sample/laboratory control sample duplicate (LCS/LCSD), matrix spike (MS) and a laboratory duplicate or matrix spike duplicate (MSD).

Blank samples are aliquots of analyte free water or Ottawa sand that are used as negative controls to verify that the sample collection, storage, preparation, and analysis system does not produce false positive results.

LCS recovery and precision are an indication of the laboratory's ability to successfully perform an analytical method in an interference-free matrix. MS recovery and precision are an indication of the laboratory's ability to successfully recover an analyte in the matrix of a specific sample or closely related sample matrices.

Laboratory duplicate LCSD, and MSD recovery and accuracy are an indication of the laboratory's ability to recover an analyte consistently. Surrogate spikes are used to evaluate accuracy, method performance, and extraction efficiency in each individual sample. The laboratory QC was within QAPP-specified limits for all parameters in the QAPP.

6.0 DISCUSSION OF RESULTS

Results from this limited Site Investigation indicate that COPCs are present in the Site soil and groundwater. The limited number of COPCs detected in both the soil and groundwater onsite suggests that the effects of each COPC to human health and the environment may be evaluated against MTCA Method A CULs for Unrestricted Land Use. For COPCs that have CUL values for both the MTCA Methods A and B, the values listed under Method A were used in this evaluation. If a COPC did not have a Method A CUL value, then the Method B value was used in this evaluation. In addition, COPC detections in groundwater were also screened against Federal and State MCLs.

Diesel- and oil-range hydrocarbons, phthalates, and metals were detected in the Site soil on the NRDS and former Oroweat properties; and tetrachloroethenePCE was also detected in one soil boring on the NRDS property. Results from this investigation indicate that oil-range hydrocarbons is the only COPC to be detected above the MTCA Method A CULs. No COPCs were detected above Method B CULs where there were no Method A values Because the Site soil does not appear to be impacted by multiple chemicals of concern, detections of the COPCs have been evaluated against MTCA Method A CULs in this section. VOCs, SVOCs, and barium were observed in the Site groundwater. Results from this investigation indicate that no COPCs were detected above MTCA Method A CULs or Federal/State MCLs. One COPC (1,1-DCE), which did not have a MTCA Method A CUL, was detected in a concentration above the MTCA Method B CUL. No other COPCs were detected above Method B CULs if there were no value for the MTCA Method A CUL.Because the Site groundwater appears to be impacted by few chemicals of concern, the groundwater COPCs have been evaluated against MTCA Method A CULs (where available) and Federal MCLs in this section.

6.1 COPCs in Site Soil

Soil samples were collected in three areas of interest onsite. One soil boring (MW-3/B-5) is located north and hydraulically upgradient of the NRDS facility operations. One soil boring (B-3) was advanced on the former Oroweat property boundary downgradient of a fuel (diesel and gasoline) and chlorinated solvent release. Two soil borings were advanced on the NRDS property downgradient of the waste transfer station facility (MW-1/B-1) and near storage buildings (MW-2/B-4). No soil samples were collected from boring B-2, located in the southern portion of the former Oroweat property.

6.1.1 Upgradient Soil Boring

One upgradient soil boring, B-5 was advanced in the landscaped area along the northern boundary of the Site (Figures 3 and 4). No staining or petroleum odors were observed in this soil boring. Oil-range hydrocarbons were detected in concentrations above MTCA Method A in the sample collected at a depth of 7 feet bgs (Figure 5). The source of the oil-range hydrocarbons is not known, and on the basis of the location of the boring and the inferred groundwater gradient onsite, the source of the oil- range hydrocarbons is suspected to have originated from an offsite source. The stormwater pipe system located north and upgradient from of the boring is a potential source.

It should be noted that tetrachloroethene PCE, a chlorinated solvent, was also detected in the 7-foot soil sample collected from boring B-5. The detected concentration of 0.005 mg/kg is below the MTCA Method A CUL of 0.05 mg/kg, and therefore this detection of tetrachloroethene PCE does not pose a human health risk (Figure 5). However, the presence of tetrachloroethene PCE indicates that a release of chlorinated solvents has occurred near the area of the soil boring or upgradient of the soil boring. The source and lateral extent of the tetrachloroethene PCE release is not known. The stormwater line north and upgradient from the boring is a potential source.

Acetone, toluene, di-n-butylphthalate, barium, and chromium were also detected in concentrations below MTCA A <u>CULs</u>. These COPCs are not anticipated to be chemicals of concern because of the following reasons: common laboratory contaminant (acetone), suspected to have originated in the laboratory container (toluene), suspected to be associated with the oil-range hydrocarbons but not at a concentration of concern (di-n-butylphthalate), and naturally occurring (barium and chromium).

6.1.2 Former Oroweat Property

One soil boring, B-3, was advanced east and downgradient of a petroleum and chlorinated solvent release (Figure 3). A faint petroleum odor was observed in the soil sample collected at 3 feet bgs. Although this soil sample was analyzed for total petroleum hydrocarbons, VOCs, and SVOCs. No COPCs was were detected in concentrations exceeding MTCA Method A CULs, and therefore, the detected concentrations of these COPCs do not pose a human health risk. The following COPCs were detected: diesel and oil range hydrocarbons, toluene, dinbutylphthalate, phenanthrene, and pyrene. It should be noted that wWith the exception of the detection of toluene, the presence of these compounds are suspected to be associated with the release of diesel compounds from LUSTs removed from the eastern area of this building. The detection of toluene is suspected to have originated in the laboratory container.

6.1.3 NRDS Property

Two soil borings were advanced on the NRDS property (B-1 and B-4) in the downgradient and downslope areas of the property (Figures 3 and 4). Low concentrations of oil-range hydrocarbons (B-4 only), acetone, toluene, di-n-butylphthalate, barium, and chromium were detected in these borings. None of these detections exceeded—were above MTCA Method A CULs, and therefore the detected concentrations of these COPCs do not pose a human health risk. The should be noted that the detection of oil-range hydrocarbons indicates that there may be due to the presence of a release of oil-range residual petroleum hydrocarbons remaining onsite after the removal of two USTs formerly located near—northeast and southwest of B-4—or upgradient from B-4 (Figure 6). The source and the lateral extent of the oil-range hydrocarbons is not known. The remaining detections are not anticipated to be chemicals of concern because of the following reasons: common laboratory contaminant (acetone), suspected to have originated in the laboratory container (toluene), suspected to be associated with the oil-range hydrocarbons but not at a concentration of concern (di-n-butylphthalate), and naturally occurring (barium and chromium).

6.2 COPCs in the Site Groundwater

Groundwater samples were collected from three monitoring wells installed on the NRDS property. One monitoring well is located upgradient of the facility (MW-3/B-5), one monitoring well is located downgradient of the facility (MW-1/B-1), and one monitoring well is located adjacent to storage buildings (MW-2/B-4). Detections of the COPCs in the groundwater are discussed by the location of the wells.

6.2.1 Upgradient Monitoring Well

MW-3/B-5 is considered the most hydraulically up-gradient well on the Site (Figure 3) on the basis of the Site topography. TetrachloroethenePCE and barium were detected in the groundwater collected from this monitoring well in concentrations below MTCA Method A CULs and federal/state MCLs. No COPCs were detected in concentrations above the MTCA Method B CULs when no value was available for the MTCA Method A CULs. TetrachloroethenePCE is a solvent, and its presence in the upgradient well indicates that this compound has originated from an offsite source and may be migrating onto the NRDS property (Figure 6). The detected concentration does not pose a human health risk; however, the source of this chlorinated solvent and the lateral extent of this compound is not known.

6.2.2 Downgradient Monitoring Wells

Two monitoring wells, MW-1/B-1 and MW-2/B-4 are located in the southern portion of the Site. 1,1,1-Trichloroethane; 1,1-dichloroethaneDCA; 1,1-dichloroethane; 1,2-dichloroethaneDCA; benzene; chloroform; dibromomethane; ethylbenzene; xylene; di-n-butylphthalate; pyrene; and barium were detected in the groundwater collected from one or both of these wells in concentrations below MTCA Method A CULs and federal/state MCLs. One COPC (1,1-DCE) did not have a MTCA Method A CUL, and the detected concentration was above the MTCA Method B CULs when no value was available for the MTCA Method A CULs.

The presence of solvents in MW-2/B-4 (Figure 6) indicates that there may <u>have been</u> a release of solvents near or upgradient of MW-2/B-4, and the presence of ethylbenzene and xylene in MW-1/B-1 indicates that there may <u>have been</u> a release of gasoline near or upgradient of MW1/B1. Potential areas include the NRDS main building and storm drain system. In addition, the solvents detected in MW-2 may be associated with a historic release from a former carpet cleaning operation on the Site which was located east and upgradient of MW-2/B-4 or from the release reported on the former Oroweat property. The carpet cleaning operation is suspected to have been present onsite between 1930 and 196536 (AMEC 2008).

7.0 CONCLUSIONS AND RECOMMENDATIONS

Results from this limited Site Investigation indicate that COPCs are present in the Site soil and groundwater MTCA A CULs. These COPCs include diesel- and oil- range hydrocarbons, VOCs, SVOCs, barium and chromium in the soil; and constituents of gasoline, solvents (VOCs), SVOCs, and barium in the groundwater. Diesel and oil range hydrocarbons were detected in one soil sample at a concentration exceeding MTCA Method A CULs, and the remaining COPCs detected in soil were below MTCA Method A CULs. None of the detected concentrations exceed MTCA Method A CULs or Federal MCLs for drinking water. Therefore, the detected concentrations of these COPCs do not pose a human health risk. However, the presence of these chemicals indicates that there may be may have been a release of these compounds near or upgradient of the soil and groundwater sample locations.

Two USTs were removed from the NRDS property in 1994, and a release of fuel was reported to Ecology. <u>Upon discovery of the petroleum-impacted soil.</u> No additional records indicating that the release was mitigated were on file at Ecology. The location of the two USTs removed on the NRDS property was not described in the documents reviewed. If SPU has an independent cleanup was conducted, and petroleum-impacted soil on the NRDS property was removed to the extent practical to maintain the structural integrity of two retaining walls and a building located in the southeast corner of the property. Observations recorded during the cleanup activities described the visible impacted soil in the southern excavation sidewall as a "2 to 3 inch (gray-stained soil lens) 2 to 3 feet wide (Herrera 1997)." Consequently, petroleum-impacted soil was left undisturbed in the southeast corner of the property. records regarding the removal of the USTs, the records should be reviewed, and the extent of the release should be evaluated.

In addition, the 15500 NNorth 34th Street property has a restrictive covenant due to gasoline and diesel-impacted soil remaining below the building footprint. If future development plans include demolition of the building or disturbance of the impacted soil, Ecology is required to be notified, and the diesel-impacted soil will be required to be managed as <u>petroleum contaminated soilapotentially hazardous waste</u>.

On the basis of the results of this investigation, it is AMEC's opinion that the following actions be conducted onsite:

- 1. If the soil below the 15500 NNorth 34th Street building will be disturbed, Ecology should be notified in accordance with the property's restrictive covenant, and a soil management plan should be prepared. The purpose of the plan is to address 1) worker safety monitoring and hazard prevention in the known contaminated areas onsite; 2) proposed plans to segregate, manage, and dispose of contaminated soil; and equipment decontamination.
- If the NRDS is to be demolished, it is recommended that additional soil and groundwater samples be collected below building foundations in areas where floor drains are located. Gasoline constituents were detected in the Site groundwater in concentrations below MTCA CULs and State MCLs. The presence of the gasoline constituents in only one of

- the three monitoring wells indicates that a localized source may be present within the main building. The presence of the gasoline constituents in the groundwater in the central area indicates that there is a release of gasoline onsite.
- 3. Should development plans include the demolition of the NRDS building located in the southeast corner of the 1350 North 34th Street property, the soil below the building and the electrical subsurface utility line connecting this building to the main building should be evaluated for the presence of petroleum, specifically gasoline and diesel. A lens of petroleum-impacted soil was reported in these areas in 1994 during the removal of two underground storage tanks. Where practical, the petroleum contaminated soil was removed. To maintain the structural integrity of two retaining walls and a building structure, the cleanup activities did not remove all of the petroleum contaminated soil onsite. Additional cleanup could be conducted in conjunction with Site redevelopment.
- 3.Locate and review all documents pertaining to the former USTs on the NRDS property and evaluate the extent of the reported petroleum release associated with the USTs. Following this evaluation, the UST area may require additional environmental investigation and remediation to achieve closure from Ecology. This could be conducted in conjunction with Site redevelopment.
- 4. Collect groundwater from the three of the existing wells during the dry season to monitor the presence of VOCs, specifically the solvents and gasoline constituents in the groundwater. The purpose of the groundwater evaluation is to confirm that the VOCs, solvents, and gasoline constituents have remained at levels below Ecology MTCA Method A CULscleanup levels.
- 5. Evaluate the extent of the elevated oil-range hydrocarbons detected along the northern property boundary.
- 6. Survey the elevations of the top of casing in the monitoring wells onsite. The groundwater gradient onsite may be calculated following completion of this survey.

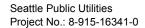
8.0_CLOSURE

AMEC appreciates the opportunity to be of service to the SPU Materials Laboratory on this project. If you have any questions or comments regarding this report, please feel free to contact us at (425) 820-4669.

AMEC Earth & Environmental, Inc.

Cherilyn Inouye Meg Strong, LG
Project Manager Senior Associate

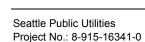
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9.08.0 LIMITATIONS

This report was prepared exclusively for the SPU Materials Laboratory by AMEC Earth & Environmental, Inc. The quality of information, conclusions, and estimates contained herein is consistent with the level of effort involved in AMEC services and based on: i) information available at the time of preparation, ii) data supplied by outside sources, and iii) the assumptions, conditions, and qualifications set forth in this report and AMEC proposal. This report is intended to be used by the SPU Materials Laboratory for the Site only, subject to the terms and conditions of the Seattle Public Utility's contract with AMEC. Any other use of, or reliance on, this report by any third party is at that party's sole risk.

The findings contained herein are relevant to the dates of the AMEC Site visit and should not be relied upon to represent conditions at later dates. Data presented herein are from discreet sampling points identified in our report, and can not be construed to represent conditions at unsampled locations.



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TABLES

Table 1

Total Petroleum Hydrocarbons Soil Analytical Results

SPU- North Recycling and Disposal Station

Seattle, Washington

| | | | | TPH-Diesel Range | TPH-Lube Oil Range | TPH-Gasoline Range |
|-------------|------------------|---------------|---------------------|------------------|--------------------|--------------------|
| | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Method | d A CULs | | | 2,000 | 2,000 | 30 |
| MTCA Method | d A CULs for Inc | dustrial Land | Use | | | |
| MTCA Method | d B CULs, Carc | inogen | | | | |
| MTCA Method | d B CULs, Non- | carcinogen | | | | |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | |
| B-1 | 3/10/2008 | B-1_5 | 5 | 29.0 U | 57.0 U | 5.60 U |
| B-3 | 3/10/2008 | B-3_3 | 3 | 150 | 300 | 6.50 U |
| B-4 | 3/11/2008 | B-4_6 | 6 | 28.0 U | 110 | 4.60 U |
| B-5 (DUP) | 3/11/2008 | B-5_6 | 7 | 140 U | 2,700 | 6.00 U |
| B-5 | 3/11/2008 | B-5_7 | 7 | 140 U | 2,400 | 5.10 U |

BOLD = detection

All results have not been not validated

mg/Kg = milligrams per kilogram

bgs = below ground surface

U = The analyte is not detected at or above the concentration listed

--- = no data/not researched

DUP = Field Duplicate

| | | | | (cis) 1,2-Dichloroethene | (cis) 1,3-Dichloropropene | (trans) 1,2-Dichloroethene | (trans) 1,3-Dichloropropene | 1,1,2-Tetrachloroethane | 1,1,1-Trichloroethane | 1,1,2,2-Tetrachloroethane | 1,1,2-Trichloroethane | 1,1-Dichloroethane | 1,1-Dichloroethene | 1,1-Dichloropropene |
|-------------|----------------------------------|-----------|------------------|--------------------------|---------------------------|----------------------------|-----------------------------|-------------------------|-----------------------|---------------------------|-----------------------|--------------------|--------------------|---------------------|
| | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Meth | | | | | | | | | 2 | | | | | |
| - | nod A CULs for | | Use | | | | | 38 | 2 | | 40 | | 4.7 | |
| | nod B CULs, Ca nod B CULs, No | | | 800 | 5.6 | 1.600 | 5.6 | | | 5 | 18 320 | 9.000 | 1.7 | |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | 000 | 2,400 | 1,000 | 2,400 | 2,400 | 72,000 | | 320 | 8,000 | 4,000 | |
| B-1 | 3/10/2008 | B-1_5 | 5 | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U |
| B-3 | 3/10/2008 | B-3_3 | 3 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U |
| B-4 | 3/11/2008 | B-4_6 | 6 | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U |
| B-5 (DUP) | 3/11/2008 | B-5_6 | 7 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U |
| B-5 | 3/11/2008 | B-5_7 | 7 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) from *Washington State Department of Ecology, Cleanup Levels* and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

| | | | 1,2,3-Trichlorobenzene | 1,2,3-Trichloropropane | 1,2,4-Trichlorobenzene | 1,2,4-Trimethylbenzene | 1,2-Dibromo-3-chloropropane | 1,2-Dibromoethane | 1,2-Dichlorobenzene | 1,2-Dichloroethane | 1,2-Dichloropropane | 1,3,5-Trimethylbenzene | 1,3-Dichlorobenzene |
|-----------|----------------|--------------|------------------------|------------------------|------------------------|------------------------|-----------------------------|-------------------|---------------------|--------------------|---------------------|------------------------|---------------------|
| | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| | nod A CULs | | | | | | | | | | | | |
| | nod A CULs for | | | | | | | | | | | | |
| | nod B CULs, Ca | | | 0.14 | | | 0.71 | | | 11 | 15 | | |
| Location | nod B CULs, No | n-carcinogen | | 480 | 800 | 4,000 | | | 7,200 | 1,600 | | 4,000 | |
| ID | Sample Date | Sample ID | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_5 | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.005 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U |
| B-3 | 3/10/2008 | B-3_3 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.005 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U |
| B-4 | 3/11/2008 | B-4_6 | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.005 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U |
| B-5 (DUP) | 3/11/2008 | B-5_6 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.006 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U |
| B-5 | 3/11/2008 | B-5_7 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.005 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface --- = no data/not researched

U = The analyte is not detected at or above the listed reporting detection limit

DUP = Field duplicate

| | | | 1,3-Dichloropropane | 1,4-Dichlorobenzene | 2,2-Dichloropropane | 2-Butanone | 2-Chloroethyl Vinyl Ether | 2-Chlorotoluene | 2-Hexanone | 4-Chlorotoluene | Acetone | Benzene | Bromobenzene |
|----------------|----------------|--------------|---------------------|---------------------|---------------------|------------|---------------------------|-----------------|------------|-----------------|---------|----------|--------------|
| | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | | (mg/Kg) |
| MTCA Meth | | | | | | | | | | | | 0.03 | |
| | nod A CULs for | | | | | | | | | | | 0.03 | |
| | nod B CULs, Ca | | | 42 | | | | | | | | 18 | |
| | od B CULs, No | n-carcinogen | | | | | | | | | 8,000 | 320 | |
| Location ID | Sample Date | Sample ID | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_5 | 0.0009 U | 0.0009 U | 0.0009 U | 0.005 U | 0.005 U | 0.0009 U | 0.005 U | 0.0009 U | 0.014 | 0.0009 U | 0.0009 U |
| B-3 | 3/10/2008 | B-3_3 | 0.001 U | 0.001 U | 0.001 U | 0.005 U | 0.005 U | 0.001 U | 0.005 U | 0.001 U | 0.005 U | 0.001 U | 0.001 U |
| B-4 | 3/11/2008 | B-4_6 | 0.0010 U | 0.0010 U | 0.0010 U | 0.005 U | 0.005 U | 0.0010 U | 0.005 U | 0.0010 U | 0.005 U | 0.0010 U | 0.0010 U |
| B-5 (DUP) | 3/11/2008 | B-5_6 | 0.001 U | 0.001 U | 0.001 U | 0.006 U | 0.006 U | 0.001 U | 0.006 U | 0.001 U | 0.730 | 0.001 U | 0.001 U |
| B-5 | 3/11/2008 | B-5_7 | 0.001 U | 0.001 U | 0.001 U | 0.005 U | 0.005 U | 0.001 U | 0.005 U | 0.001 U | 0.380 | 0.001 U | 0.001 U |

Notes:

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U = The analyte is not detected at or above the listed reporting detection limit

DUP = Field duplicate

| | | | Bromochloromethane | Bromodichloromethane | Bromoform | Bromomethane | Carbon Disulfide | Carbon Tetrachloride | Chlorobenzene | Chloroethane | Chloroform | Chloromethane | Dibromochloromethane |
|----------------|----------------|--------------|--------------------|----------------------|-----------|--------------|------------------|----------------------|---------------|--------------|------------|---------------|----------------------|
| | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Meth | | | | | | | | | | | | | |
| | nod A CULs for | | | | | | | | | | | | |
| | nod B CULs, Ca | | | 16 | 130 | | | 7.7 | | | 160 | 77 | 12 |
| | nod B CULs, No | n-carcinogen | | 1,600 | 1,600 | 110 | 8,000 | 56 | 70,000 | 1,600 | 800 | | 1,600 |
| Location ID | Sample Date | Sample ID | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_5 | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.005 U | 0.0009 U | 0.005 U | 0.0009 U |
| B-3 | 3/10/2008 | B-3_3 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.005 U | 0.001 U | 0.005 U | 0.001 U |
| B-4 | 3/11/2008 | B-4_6 | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.005 U | 0.0010 U | 0.005 U | 0.0010 U |
| B-5 (DUP) | 3/11/2008 | B-5_6 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.006 U | 0.001 U | 0.006 U | 0.001 U |
| B-5 | 3/11/2008 | B-5_7 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.005 U | 0.001 U | 0.005 U | 0.001 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface --- = no data/not researched

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DUP = Field duplicate

| | | | Dibromomethane | Dichlorodifluoromethane | Ethylbenzene | Hexachlorobutadiene | lodomethane | Isopropylbenzene | Methyl Isobutyl Ketone | Methyl t-Butyl Ether | Methylene Chloride | Naphthalene | n-Butylbenzene |
|-----------|----------------|--------------|----------------|-------------------------|--------------|---------------------|-------------|------------------|------------------------|----------------------|--------------------|-------------|----------------|
| | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Meth | | | | | 6 | | | | | 0.1 | 0.02 | 5 | |
| | nod A CULs for | | | | 6 | | | | | 0.1 | 0.02 | 5 | |
| | od B CULs, Ca | - | | 40.000 | | 13 | | | | 560 | 130 | 4.000 | |
| Location | od B CULs, No | n-carcinogen | | 16,000 | 8,000 | 16 | | | 6,400 | 69,000 | 4,800 | 1,600 | |
| ID | Sample Date | Sample ID | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_5 | 0.0009 U | 0.0009 U | 0.0009 U | 0.005 U | 0.005 U | 0.0009 U | 0.005 U | 0.0009 U | 0.005 U | 0.0009 U | 0.0009 U |
| B-3 | 3/10/2008 | B-3_3 | 0.001 U | 0.001 U | 0.001 U | 0.005 U | 0.005 U | 0.001 U | 0.005 U | 0.001 U | 0.005 U | 0.001 U | 0.001 U |
| B-4 | 3/11/2008 | B-4_6 | 0.0010 U | 0.0010 U | 0.0010 U | 0.005 U | 0.005 U | 0.0010 U | 0.005 U | 0.0010 U | 0.005 U | 0.0010 U | 0.0010 U |
| B-5 (DUP) | 3/11/2008 | B-5_6 | 0.001 U | 0.001 U | 0.001 U | 0.006 U | 0.006 U | 0.001 U | 0.006 U | 0.001 U | 0.006 U | 0.001 U | 0.001 U |
| B-5 | 3/11/2008 | B-5_7 | 0.001 U | 0.001 U | 0.001 U | 0.005 U | 0.005 U | 0.001 U | 0.005 U | 0.001 U | 0.005 U | 0.001 U | 0.001 U |

Notes:

BOLD = detection

DUP = Field duplicate

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface --- = no data/not researched

U = The analyte is not detected at or above the listed reporting detection limit

| | | | n-Propylbenzene | p-IsopropyItoluene | sec-Butylbenzene | Styrene | tert-Butylbenzene | Tetrachloroethene | Toluene | Trichloroethene | Trichlorofluoromethane | Vinyl Acetate | Vinyl Chloride |
|-----------|----------------|--------------|-----------------|--------------------|------------------|----------|-------------------|-------------------|---------|-----------------|------------------------|---------------|----------------|
| | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Meth | | | | | | | | 0.05 | 7 | 0.03 | | | |
| | nod A CULs for | | | | | | | 0.05 | 7 | 0.03 | | | 0.07 |
| | nod B CULs, Ca | | | | | 33 | | 240 35,000 | 290,000 | 1.9 800 | 24.000 | 90,000 | 0.67 |
| Location | nod B CULs, No | n-carcinogen | | | | 16,000 | | 35,000 | 280,000 | 600 | 24,000 | 80,000 | 240 |
| ID | Sample Date | Sample ID | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_5 | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.0009 U | 0.002 | 0.0009 U | 0.0009 U | 0.005 U | 0.0009 U |
| B-3 | 3/10/2008 | B-3_3 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.003 | 0.001 U | 0.001 U | 0.005 U | 0.001 U |
| B-4 | 3/11/2008 | B-4_6 | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.0010 U | 0.003 | 0.0010 U | 0.0010 U | 0.005 U | 0.0010 U |
| B-5 (DUP) | 3/11/2008 | B-5_6 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.005 | 0.002 | 0.001 U | 0.001 U | 0.006 U | 0.001 U |
| B-5 | 3/11/2008 | B-5_7 | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 U | 0.001 | 0.003 | 0.001 U | 0.001 U | 0.005 U | 0.001 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) from *Washington State Department of Ecology, Cleanup Levels* and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

| | | | m,p-Xylene | o-Xylene | Total-Xylene |
|----------------|----------------|-----------------|------------|----------|--------------|
| | | | (mg/Kg) | (mg/Kg) | (mg/kg) |
| | nod A CULs | | | | 9 |
| MTCA Meth | nod A CULs for | Industrial Land | | | 9 |
| MTCA Meth | nod B CULs, Ca | ırcinogen | | | |
| | nod B CULs, No | n-carcinogen | 160,000 | 160,000 | 16,000 |
| Location ID | Sample Date | Sample ID | | | |
| B-1 | 3/10/2008 | B-1_5 | 0.002 U | 0.0009 U | 0.002 |
| B-3 | 3/10/2008 | B-3_3 | 0.002 U | 0.001 U | 0.002 |
| B-4 | 3/11/2008 | B-4_6 | 0.002 U | 0.0010 U | 0.002 |
| B-5 (DUP) | 3/11/2008 | B-5_6 | 0.002 U | 0.001 U | 0.002 |
| B-5 | 3/11/2008 | B-5_7 | 0.002 U | 0.001 U | 0.002 |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface --- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) from *Washington State Department of Ecology, Cleanup Levels* and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or above the listed reporting detection limit

| | | | | (3+4)-Methylphenol (m,p-Cresol) | 1,2,4-Trichlorobenzene | 1,2-Dichlorobenzene | 1,2-Dinitrobenzene | 1,2-Diphenylhydrazine | 1,3-Dichlorobenzene | 1,3-Dinitrobenzene | 1,4-Dichlorobenzene | 1,4-Dinitrobenzene | 1-Methylnaphthalene | 2,3,4,6-Tetrachlorophenol |
|-------------|---|------------|---------------------|------------------------------------|------------------------|---------------------|--------------------|-----------------------|---------------------|--------------------|---------------------|--------------------|---------------------|---------------------------|
| | 24 Mathad 4 CIII s | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| - | CA Method A CULs | | | | | | | | | | | | | |
| | CA Method A CULs CA Method A CULs for Industrial Land Use | | | | | | | | | | | | | |
| MTCA Method | B CULs, Carc | inogen | | | | | | 1.3 | | | 42 | | | |
| MTCA Method | B CULs, Non- | carcinogen | | | 800 | 7,200 | | | | | | | 24 | 2,400 |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_8 | 8 | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.007 U | 0.037 U |
| B-3 | 3/10/2008 | B-3_3 | 3 | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.008 U | 0.041 U |
| B-4 | 3/11/2008 | B-4_8 | 8 | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.007 U | 0.037 U |
| B-5 (DUP) | 3/11/2008 | B-5_8 | 9 | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.007 U | 0.035 U |
| B-5 | 3/11/2008 | B-5_9 | 9 | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.007 U | 0.035 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) from *Washington State Department of Ecology,* Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or above the listed reporting detection limit

| | | | | (S) 2,3,5,6-Tetrachlorophenol | 2,3-Dichloroaniline | 2,4,5-Trichlorophenol | 2,4,6-Trichlorophenol | 2,4-Dichlorophenol | 2,4-Dimethylphenol | 2,4-Dinitrophenol | 2,4-Dinitrotoluene | 2,6-Dinitrotoluene | 2-Chloronaphthalene | 2-Chlorophenol |
|-------------|---|-------|---------------------|-------------------------------|---------------------|-----------------------|-----------------------|--------------------|--------------------|-------------------|--------------------|--------------------|---------------------|----------------|
| MTCA Mada | CA Method A CULs | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | | | | (mg/Kg) | | | (mg/Kg) |
| | CA Method A CULs | | | | | | | | | | | | | |
| | CA Method A CULs for Industrial Land Use CA Method B CULs, Carcinogen | | | | | | 91 | | | | | | | |
| | d B CULs, Non- | | | | | 8,000 | | 240 | 1,600 | 160 | 160 | 80 | | 40 |
| Location ID | Sample Date | | Depth (feet bgs) | | | 3,000 | | 210 | 1,000 | 100 | | | | 10 |
| B-1 | 3/10/2008 | B-1_8 | 8 | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.190 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U |
| B-3 | 3/10/2008 | B-3_3 | 3 | 0.041 U | 0.041 U | 0.041 U | | | 0.041 U | | | | | |
| B-4 | 3/11/2008 | B-4_8 | 8 | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.180 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U |
| B-5 (DUP) | 3/11/2008 | B-5_8 | 9 | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.170 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U |
| B-5 | 3/11/2008 | B-5_9 | 9 | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.170 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) from Washington State Department of Ecology,

| | | | | 2-Methylnaphthalene | 2-Methylphenol (o-Cresol) | 2-Nitroaniline | 2-Nitrophenol | 3,3'-Dichlorobenzidine | 3-Nitroaniline | 4,6-Dinitro-2-methylphenol | 4-Bromophenyl-phenylether | 4-Chloro-3-methylphenol | 4-Chloroaniline | 4-Chlorophenyl-phenylether |
|-------------|--|------------|---------------------|---------------------|---------------------------|----------------|---------------|------------------------|----------------|----------------------------|---------------------------|-------------------------|-----------------|----------------------------|
| | A Method A CUI s | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| | CA Method A CULs | | | | | | | | | | | | | |
| MTCA Method | CA Method A CULs CA Method A CULs for Industrial Land Use | | | | | | | | | | | | | |
| MTCA Method | d B CULs, Carc | inogen | | | | | | 2.2 | | | | | | |
| MTCA Method | d B CULs, Non- | carcinogen | | 320 | | | | | | | | | | |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_8 | 8 | 0.007 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.190 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U |
| B-3 | 3/10/2008 | B-3_3 | 3 | 0.008 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.210 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U |
| B-4 | 3/11/2008 | B-4_8 | 8 | 0.007 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.180 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U |
| B-5 (DUP) | 3/11/2008 | B-5_8 | 9 | 0.007 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.170 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U |
| B-5 | 3/11/2008 | B-5_9 | 9 | 0.007 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.170 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) from *Washington State Department of Ecology,* Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

| | | | | 4-Nitroaniline | 4-Nitrophenol | Acenaphthene | Acenaphthylene | Aniline | Anthracene | Benzidine | Benzo[g,h,i]perylene | Benzyl alcohol | bis(2-Chloroethoxy)methane | bis(2-Chloroethyl)ether |
|-------------|--|------------|---------------------|----------------|---------------|--------------|----------------|---------|------------|-----------|----------------------|----------------|----------------------------|-------------------------|
| | CA Method A CIII s | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| | CA Method A CULs | | | | | | | | | | | | | |
| | CA Method A CULs for Industrial Land Use | | | | | | | | | | | | | |
| MTCA Method | d B CULs, Card | inogen | | | | | | 180 | | 4,300 | | | | 0.91 |
| MTCA Method | d B CULs, Non- | carcinogen | | | | 4,800 | | | 24,000 | 240 | | 24,000 | | |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_8 | 8 | 0.037 U | 0.037 U | 0.007 U | 0.007 U | 0.037 U | 0.007 U | 0.370 U | 0.007 U | 0.037 U | 0.037 U | 0.037 U |
| B-3 | 3/10/2008 | B-3_3 | 3 | 0.041 U | 0.041 U | 0.008 U | 0.008 U | 0.041 U | 0.008 U | 0.410 U | 0.008 U | 0.041 U | 0.041 U | 0.041 U |
| B-4 | 3/11/2008 | B-4_8 | 8 | 0.037 U | 0.037 U | 0.007 U | 0.007 U | 0.037 U | 0.007 U | 0.370 U | 0.007 U | 0.037 U | 0.037 U | 0.037 U |
| B-5 (DUP) | 3/11/2008 | B-5_8 | 9 | 0.035 U | 0.035 U | 0.007 U | 0.007 U | 0.035 U | 0.007 U | 0.350 U | 0.007 U | 0.035 U | 0.035 U | 0.035 U |
| B-5 | 3/11/2008 | B-5_9 | 9 | 0.035 U | 0.035 U | 0.007 U | 0.007 U | 0.035 U | 0.007 U | 0.350 U | 0.007 U | 0.035 U | 0.035 U | 0.035 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

DUP = Field duplicate

| | | | | bis(2-Chloroisopropyl)ether | bis(2-Ethylhexyl)phthalate | bis-2-Ethylhexyladipate | Butylbenzylphthalate | Carbazole | Dibenzofuran | Diethylphthalate | Dimethylphthalate | Di-n-butylphthalate | Di-n-octylphthalate | Fluoranthene |
|-------------|------------------|---------------|---------------------|-----------------------------|----------------------------|-------------------------|----------------------|-----------|--------------|------------------|-------------------|---------------------|---------------------|--------------|
| | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Method | d A CULs | | | | | | | | | | | | | |
| MTCA Method | d A CULs for Inc | dustrial Land | Use | | | | | | | | | | | |
| MTCA Method | d B CULs, Carc | inogen | | | 71 | | | 50 | | | | | | |
| MTCA Method | d B CULs, Non- | carcinogen | | 3,200 | 1,600 | | 16,000 | | 160 | 64,000 | 8,000 | 8,000 | 1,600 | 3,200 |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_8 | 8 | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.041 | 0.037 U | 0.007 U |
| B-3 | 3/10/2008 | B-3_3 | 3 | 0.041 U | 0.052 | 0.041 U | | | 0.041 U | | 0.041 U | 0.071 | 0.041 U | 0.008 U |
| B-4 | 3/11/2008 | B-4_8 | 8 | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.007 U |
| B-5 (DUP) | 3/11/2008 | B-5_8 | 9 | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.007 U |
| B-5 | 3/11/2008 | B-5_9 | 9 | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.037 | 0.035 U | 0.007 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) from *Washington State Department of Ecology,* Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

| | | | | Fluorene | Hexachlorobenzene | Hexachlorobutadiene | Hexachlorocyclopentadiene | Hexachloroethane | Isophorone | Naphthalene | Nitrobenzene | N-Nitrosodimethylamine | N-Nitroso-di-n-propylamine | N-Nitrosodiphenylamine |
|-------------|------------------|---------------|---------------------|----------|-------------------|---------------------|---------------------------|------------------|------------|-------------|--------------|------------------------|----------------------------|------------------------|
| | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Method | d A CULs | | | | | | | | | 5 | | | | |
| MTCA Method | d A CULs for Inc | dustrial Land | Use | | | | | | | 5 | | | | |
| MTCA Method | d B CULs, Carc | inogen | | | 0.63 | 13 | 0.56 | 71 | 1,100 | | | 0.02 | 0.14 | |
| MTCA Method | d B CULs, Non- | carcinogen | | 3,200 | 64 | 16 | | 80 | 16,000 | 1,600 | 40 | | | |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_8 | 8 | 0.007 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.007 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U |
| B-3 | 3/10/2008 | B-3_3 | 3 | 0.008 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U | 0.008 U | 0.041 U | 0.041 U | 0.041 U | 0.041 U |
| B-4 | 3/11/2008 | B-4_8 | 8 | 0.007 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U | 0.007 U | 0.037 U | 0.037 U | 0.037 U | 0.037 U |
| B-5 (DUP) | 3/11/2008 | B-5_8 | 9 | 0.007 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.007 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U |
| B-5 | 3/11/2008 | B-5_9 | 9 | 0.007 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U | 0.007 U | 0.035 U | 0.035 U | 0.035 U | 0.035 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched DUP = Field duplicate

 ${\tt MTCA~Method~A~and~B~Cleanup~Levels~(CULs)~from~\textit{Washington~State~Department~of~Ecology},}$

| | | | | Pentachlorophenol | Phenanthrene | Phenol | Pyrene | Pyridine | Benzo[a]anthracene | Benzo[a]pyrene | Benzo[b]fluoranthene | Benzo[k]fluoranthene | Chrysene | Dibenz[a,h]anthracene | Indeno[1,2,3-cd]pyrene | Total TEF |
|-------------|----------------|---------------|---------------------|-------------------|--------------|---------|---------|----------|--------------------|----------------|----------------------|----------------------|----------|-----------------------|------------------------|-----------|
| | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Method | A CULs | | | | | | | | | 0.1 | | | | | | 0.1 |
| MTCA Method | A CULs for Inc | dustrial Land | d Use | | | | | | | 2 | | | | | | 2 |
| MTCA Method | B CULs, Carc | inogen | | 83 | | | | | | 0.14 | | | | | | |
| MTCA Method | B CULs, Non- | carcinogen | | 2400 | | 48,000 | 2,400 | 80 | | | | | | | | |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_8 | 8 | 0.190 U | 0.007 U | 0.037 U | 0.007 U | 0.037 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.005 |
| B-3 | 3/10/2008 | B-3_3 | 3 | 0.210 U | 0.013 | 0.041 U | 0.011 | 0.041 U | 0.008 U | 0.008 U | 0.008 U | 0.008 U | 0.008 U | 0.008 U | 0.008 U | 0.006 |
| B-4 | 3/11/2008 | B-4_8 | 8 | 0.180 U | 0.007 U | 0.037 U | 0.007 U | 0.037 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.005 |
| B-5 (DUP) | 3/11/2008 | B-5_8 | 9 | 0.170 U | 0.007 U | 0.035 U | 0.007 U | 0.035 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.005 |
| B-5 | 3/11/2008 | B-5_9 | 9 | 0.170 U | 0.007 U | 0.035 U | 0.007 U | 0.035 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.007 U | 0.005 |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

--- = no data/not researched
DUP = Field duplicate

ed

U = The analyte is not detected at or above the listed reporting detection limit

TEF = Toxicity Equivalency Facto, calculated in accordance

with WAC 173-340-708(e)

MTCA Method A and B Cleanup Levels (CULs) from Washington State Department of Ecology,

Table 4 Metal Soil Analytical Results SPU- North Recycling and Disposal Station Seattle, Washington

| | | | | Arsenic | Barium | Cadmium | Total Chromium | Soluble Hexavalent Chromium | Lead | Mercury | Selenium | Silver |
|-------------|---------------|--------------|---------------------|---------|---------|---------|----------------|-----------------------------------|---------|---------|----------|---------|
| | | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Meth | od A CULs | | | 20 | | 2 | 2,000* | 19 | 250 | 2 | | |
| MTCA Meth | od A CULs for | nd Use | 20 | | 2 | 2,000 | 19 | 250 | 2 | | | |
| MTCA Meth | od B CULs, Ca | rcinogen | | 0.67 | | | | | | | | |
| MTCA Meth | od B CULs, No | n-carcinoger | ì | 24 | 16,000 | 80 | 120,000* | | 24 | 24 | 400 | 400 |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_3 | 3 | 11.0 U | 46.0 | 0.570 U | 32.0 | 1.10 U | 5.70 U | 0.280 U | 11.0 U | 0.570 U |
| B-3 | 3/10/2008 | B-3_6 | 6 | 11.0 U | 46.0 | 0.560 U | 22.0 | 1.10 U | 5.60 U | 0.280 U | 11.0 U | 0.560 U |
| B-4 | 3/11/2008 | B-4_8 | 8 | 11.0 U | 33.0 | 0.550 U | 18.0 | 1.10 U | 5.50 U | 0.270 U | 11.0 U | 0.550 U |
| B-5 (DUP) | 3/11/2008 | B-5_11 | 12 | 11.0 U | 34.0 | 0.570 U | 26.0 | 1.10 U | 5.70 U | 0.290 U | 11.0 U | 0.570 U |
| B-5 | 3/11/2008 | B-5_12 | 12 | 11.0 U | 39.0 | 0.570 U | 30.0 | 1.10 U | 5.70 U | 0.280 U | 11.0 U | 0.570 U |

Notes:

BOLD = detection

All results have not been not validated

mg/Kg = milligrams per kilogram bgs = below ground surface

--- = no data/not researched

U = The analyte is not detected at or above the listed reporting detection limit

*=Trivalent chromium CUL

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) from Washington State Department of Ecology,

Table 5 Pesticide Soil Analytical Results SPU- North Recycling and Disposal Station Seattle, Washington

| | | | | 4,4'-DDD | 4,4'-DDE | 4,4'-DDT | Aldrin | alpha-BHC | alpha-Chlordane | beta-BHC | delta-BHC | Dieldrin | Endosulfan I | Endosulfan II |
|-------------|-----------------|---------------|---------------------|----------|----------|----------|----------|-----------|-----------------|----------|-----------|----------|--------------|---------------|
| | M. J. A. O.I.II | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Met | A Method A CULs | | | | | 3 | | | | | | | | |
| MTCA Met | hod A CULs for | Industrial La | and Use | | | 4 | | | | | | | | |
| MTCA Met | hod B CULs, Ca | arcinogen | | 4 | 2.9 | 2.9 | 0.059 | | | | | 0.063 | | |
| MTCA Met | hod B CULs, No | on-carcinoge | n | | | 40 | 2 | | | | | 4 | | |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_3 | 3 | 0.011 U | 0.011 U | 0.011 U | 0.0057 U | 0.0057 U | 0.011 U | 0.0057 U | 0.0057 U | 0.011 U | 0.0057 U | 0.011 U |
| B-3 | 3/10/2008 | B-3_6 | 6 | 0.01 U | 0.01 U | 0.01 U | 0.005 U | 0.005 U | 0.01 U | 0.005 U | 0.005 U | 0.01 U | 0.005 U | 0.01 U |
| B-4 | 3/11/2008 | B-4_3 | 3 | 0.011 U | 0.011 U | 0.011 U | 0.0056 U | 0.0056 U | 0.011 U | 0.0056 U | 0.0056 U | 0.011 U | 0.0056 U | 0.011 U |
| B-5 | 3/11/2008 | B-5_3 | 3 | 0.011 U | 0.011 U | 0.011 U | 0.0054 U | 0.0054 U | 0.011 U | 0.0054 U | 0.0054 U | 0.011 U | 0.0054 U | 0.011 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

Table 5 Pesticide Soil Analytical Results SPU- North Recycling and Disposal Station Seattle, Washington

| | | | Endrin | Endrin Aldehyde | Endrin Ketone | Endsulfan Sulfate | gamma-BHC | gamma-Chlordane | Heptachlor | Heptachlor Epoxide | Methoxychlor | Toxaphene |
|-------------|-------------------|---------------|---------|-----------------|---------------|-------------------|-----------|-----------------|------------|--------------------|--------------|-----------|
| | A Method A CUII s | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) |
| MTCA Met | CA Method A CULs | | | | | | | | | | | |
| MTCA Met | nod A CULs for | Industrial La | | | | | | | | | | |
| MTCA Meti | nod B CULs, Ca | arcinogen | | | | | | | 0.22 | 0.11 | | |
| MTCA Meti | nod B CULs, No | on-carcinoge | 24 | | | | | | 40 | 1 | 400 | |
| Location ID | Sample Date | Sample ID | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_3 | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.0057 U | 0.011 U | 0.0057 U | 0.0057 U | 0.011 U | 0.0057 U |
| B-3 | 3/10/2008 | B-3_6 | 0.01 U | 0.01 U | 0.01 U | 0.01 U | 0.005 U | 0.01 U | 0.005 U | 0.005 U | 0.01 U | 0.005 U |
| B-4 | 3/11/2008 | B-4_3 | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.0056 U | 0.011 U | 0.0056 U | 0.0056 U | 0.011 U | 0.0056 U |
| B-5 | 3/11/2008 | B-5_3 | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.0054 U | 0.011 U | 0.0054 U | 0.0054 U | 0.011 U | 0.0054 U |

Notes:

BOLD = detection

All results have not been not validated mg/kg = milligrams per kilogram bgs = below ground surface

U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

Table 6 Herbicide Soil Analytical Results SPU- North Recycling and Disposal Station Seattle, Washington

| | | | | (mg/kg) 2,4,5-T | 2,4,5-TP (Silvex) | 2,4-D | 2,4-DB | Dalapon | Dicamba | Dichlorprop | Dinoseb | MCPA | MCPP | . Pentachlorophenol |
|-------------|------------------|--------------|---------------------|--------------------|-------------------|---------|---------|---------|---------|-------------|---------|---------|---------|---------------------|
| NATOA NA 41 | CA Method A CULs | | | | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) | (mg/kg) |
| | CA Method A CULs | | | | | | | | | | | | | |
| MTCA Meth | od A CULs for | nd Use | | | | | | | | | | | | |
| MTCA Meth | od B CULs, Ca | rcinogen | | | | | | | | | | | | 8 |
| MTCA Meth | od B CULs, No | n-carcinoger | ſ | 800 | 650 | 800 | 640 | 24,000 | 24,000 | | 80 | | | 24,000 |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_3 | 3 | 0.054 U | 0.054 U | 0.053 U | 0.054 U | 0.26 U | 0.053 U | 0.054 U | 0.054 U | 5.3 U | 5.3 U | 0.0011 U |
| B-3 | 3/10/2008 | B-3_6 | 6 | 0.053 U | 0.053 U | 0.052 U | 0.053 U | 0.25 U | 0.052 U | 0.052 U | 0.053 U | 5.2 U | 5.2 U | 0.0011 U |
| B-4 | 3/11/2008 | B-4_3 | 3 | 0.053 U | 0.053 U | 0.053 U | 0.053 U | 0.26 U | 0.053 U | 0.053 U | 0.053 U | 5.3 U | 5.3 U | 0.0011 U |
| B-5 | 3/11/2008 | B-5_3 | 3 | 0.051 U | 0.051 U | 0.051 U | 0.051 U | 0.25 U | 0.051 U | 0.051 U | 0.051 U | 5 U | 5 U | 0.001 U |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

Table 7 Polychlorinated Biphenyl Soil Analytical Results SPU- North Recycling and Disposal Station Seattle, Washington

| | | | | Aroclor 1016 | Aroclor 1221 | Aroclor 1232 | Aroclor 1242 | Aroclor 1248 | Aroclor 1254 | Aroclor 1260 | Aroclor 1262 | Aroclor 1268 | Total Aroclors |
|-------------|---|------------|---------------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|----------------|
| | | | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | (mg/Kg) | |
| MTCA Meth | TCA Method A CULs | | | | | | | | | | | | 1 |
| MTCA Meth | TCA Method A CULs for Industrial Land Use | | | | | | | | | | | | 10 |
| MTCA Meth | FCA Method A CULs for Industrial Land Use FCA Method B CULs, Carcinogen | | | | | | | | | | | | 0.5 |
| MTCA Meth | nod B CULs, Non | carcinogen | | | | | | | | | | | |
| Location ID | Sample Date | Sample ID | Depth (feet bgs) | | | | | | | | | | |
| B-1 | 3/10/2008 | B-1_3 | 3 | 0.057 U | ND |
| B-3 | 3/10/2008 | B-3_6 | 6 | 0.056 U | ND |
| B-4 | 3/11/2008 | B-4_3 | 3 | 0.056 U | ND |
| B-5 | 3/11/2008 | B-5_3 | 3 | 0.054 U | ND |

Notes:

BOLD = detection

All results have not been not validated mg/Kg = milligrams per kilogram bgs = below ground surface

--- = no data/not researched

U = The analyte is not detected at or above the listed reporting detection limit

Table 8

Total Petroleum Hydrocarbon Groundwater Analytical Results
SPU- North Recycling and Disposal Station
Seattle, Washington

| | | | TPH-Diesel Range | TPH- Lube Oil Range | TPH-Gasoline Range |
|------------------|--------------------|-----------|------------------|---------------------|--------------------|
| | | | (mg/L) | (mg/L) | (µg/L) |
| MTCA Method A CL | JL | | 0.5 | 0.5 | 800 |
| Groundwater ARAR | - Federal MCL | | | | |
| MTCA Method B Cl | JLs, Carcinogen | | | | |
| MTCA Method B CL | JLs, Noncarcinogen | | | | |
| Location ID | Sample Date | Sample ID | | | |
| MW1 | 3/14/2008 | DUP-1 | 0.280 U | 0.450 U | 100 U |
| MW1 | 3/14/2008 | MW-1 | 0.270 U | 0.430 U | 100 U |
| MW2 | 3/14/2008 | MW-2 | 0.260 U | 0.420 U | 100 U |
| MW3 | 3/14/2008 | MW-3 | 0.270 U | 0.430 U | 100 U |

All results have not been not validated

TPH = total petroleum hydrocarbons

mg/L = milligrams per liter

μg/L = micrograms per liter

BOLD = detection

U = The analyte is not detected at or above the concentration listed

DUP = Field Duplicate

--- = no data/not researched

MTCA Method A Cleanup Levels (CULs) from Washington State Department of Ecology,

Table 9
Volatile Organic Compound Groundwater Analytical Results
SPU- North Recycling and Disposal Station
Seattle, Washington

| | | | (cis) 1,2-Dichloroethene | (cis) 1,3-Dichloropropene | (trans) 1,2-Dichloroethene | (trans) 1,3-Dichloropropene | 1,1,1,2-Tetrachloroethane | 1,1,1-Trichloroethane | 1,1,2,2-Tetrachloroethane | 1,1,2-Trichloroethane | 1,1-Dichloroethane | 1,1-Dichloroethene | 1,1-Dichloropropene | 1,2,3-Trichlorobenzene | 1,2,3-Trichloropropane | 1,2,4-Trichlorobenzene |
|----------------|-----------------|-----------|--------------------------|---------------------------|----------------------------|-----------------------------|---------------------------|-----------------------|---------------------------|-----------------------|--------------------|--------------------|---------------------|------------------------|------------------------|------------------------|
| | | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method A | | | | | | | | 200 | | | | | | | | |
| Groundwater AR | AR- Federal MCL | = | | | | | | 200 | | 5 | | | | | | |
| MTCA Method B | CULs, Carcinoge | en | | 0.24 | | 0.24 | 1.7 | | 0.22 | 0.77 | | 0.073 | | | 0.0063 | |
| MTCA Method B | CULs, Noncarcin | ogen | 80 | 240 | 160 | 240 | 240 | 7,200 | | 32 | 800 | 400 | | | 48 | 80 |
| Location ID | Sample Date | Sample ID | | · | | · | | | | · | | | | · | · | |
| MW1 | 3/14/2008 | DUP-1 | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U |
| MW1 | 3/14/2008 | MW-1 | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U |
| MW2 | 3/14/2008 | MW-2 | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 30.0 | 0.200 U | 0.200 U | 30.0 | 31.0 | 0.200 U | 0.200 U | 0.200 U | 0.200 U |
| MW3 | 3/14/2008 | MW-3 | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U |

BOLD = detection

All results have not been not validated µg/L = micrograms per liter U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

DUP = Field duplicate

Table 9
Volatile Organic Compound Groundwater Analytical Results
SPU- North Recycling and Disposal Station
Seattle, Washington

| | | | 1,2,4-Trimethylbenzene | 1,2-Dibromo-3-chloropropane | 1,2-Dibromoethane | 1,2-Dichlorobenzene | 1,2-Dichloroethane | 1,2-Dichloropropane | 1,3,5-Trimethylbenzene | 1,3-Dichlorobenzene | 1,3-Dichloropropane | 1,4-Dichlorobenzene | 2,2-Dichloropropane | 2-Butanone | 2-Chloroethyl Vinyl Ether | 2-Chlorotoluene |
|---------------|-----------------|-----------|------------------------|-----------------------------|-------------------|---------------------|--------------------|---------------------|------------------------|---------------------|---------------------|---------------------|---------------------|------------|---------------------------|-----------------|
| | | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method A | | | | | | | 5 | | | | | | | | | |
| | AR- Federal MCL | | | 0.4 | | 600 | 5 | 5 | | | | 75 | | | | |
| MTCA Method B | | | | 0.031 | | | 0.48 | 0.64 | | | | 1.8 | | | | |
| MTCA Method B | | | 400 | | | 720 | 160 | | 400 | | | | | | | |
| Location ID | Sample Date | Sample ID | | | | | | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 0.200 U | 1.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 5.00 U | 1.00 U | 0.200 U |
| MW1 | 3/14/2008 | MW-1 | 0.200 U | 1.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 5.00 U | 1.00 U | 0.200 U |
| MW2 | 3/14/2008 | MW-2 | 0.200 U | 1.00 U | 0.200 U | 0.200 U | 3.40 | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 5.00 U | 1.00 U | 0.200 U |
| MW3 | 3/14/2008 | MW-3 | 0.200 U | 1.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 5.00 U | 1.00 U | 0.200 U |

BOLD = detection

All results have not been not validated µg/L = micrograms per liter U = The analyte is not detected at or above the listed reporting detection limit

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DUP = Field duplicate

Table 9
Volatile Organic Compound Groundwater Analytical Results
SPU- North Recycling and Disposal Station
Seattle, Washington

| | | | 2-Hexanone | 4-Chlorotoluene | Acetone | Benzene | Bromobenzene | Bromochloromethane | Bromodichloromethane | Bromoform | Bromomethane | Carbon Disulfide | Carbon Tetrachloride | Chlorobenzene | Chloroethane | Chloroform |
|---------------------|-----------------------------|-----------|------------|-----------------|---------|---------|--------------|--------------------|----------------------|-----------|--------------|------------------|----------------------|---------------|--------------|------------|
| NATOA NA sils sid A | O. II | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method A | | | | | | 5 5 | | 80 | | 80 | | | 5 | 100 | | 80 |
| | AR- Federal MCL | | | | | 0.8 | | | 0.71 | 5.5 | | | | | | 7.2 |
| | CULs, Carcinoge | | | | 900 | 32 | | | 0.71 160 | | 11 | 900 | 0.34 | 160 | | |
| Location ID | CULs, Noncarcin Sample Date | Sample ID | | | 800 | 32 | | | 160 | 160 | 11 | 800 | 5.6 | 160 | | 80 |
| MW1 | 3/14/2008 | DUP-1 | 2.00 U | 0.200 U | 5.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U |
| MW1 | 3/14/2008 | MW-1 | 2.00 U | 0.200 U | 5.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U |
| MW2 | 3/14/2008 | MW-2 | 2.00 U | 0.200 U | 5.00 U | 0.320 | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.660 |
| MW3 | 3/14/2008 | MW-3 | 2.00 U | 0.200 U | 5.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U |

BOLD = detection

All results have not been not validated µg/L = micrograms per liter U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

DUP = Field duplicate

Table 9
Volatile Organic Compound Groundwater Analytical Results
SPU- North Recycling and Disposal Station
Seattle, Washington

| | | | Chloromethane | Dibromochloromethane | Dibromomethane | Dichlorodifluoromethane | Ethylbenzene | Hexachlorobutadiene | lodomethane | Isopropylbenzene | Methyl Isobutyl Ketone | Methyl t-Butyl Ether | Methylene Chloride | Naphthalene | n-Butylbenzene | n-Propylbenzene |
|------------------|-----------------|-----------|---------------|----------------------|----------------|-------------------------|--------------|---------------------|-------------|------------------|------------------------|----------------------|--------------------|-------------|----------------|-----------------|
| NATION NA (L. LA | 0.11 | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method A | | | | | | | 700 | | | | | 20 | 5 | 160 | | |
| | AR- Federal MCL | | | 80 | | | 700 | 0.50 | | | | | 5 | | | |
| MTCA Method B | | | 3.4 | 0.52 | | 4 000 | | 0.56 | | | 0.40 | 24 | 5.8 | 400 | | |
| MTCA Method B | · | | | 160 | | 1,600 | 800 | 1.6 | | | 640 | 6,900 | 480 | 160 | | |
| Location ID | Sample Date | Sample ID | | | | | | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 1.00 U | 0.310 | 0.200 U | 0.200 U | 0.280 | 0.200 U | 1.00 U | 0.200 U | 2.00 U | 0.200 U | 1.00 U | 1.00 U | 0.200 U | 0.200 U |
| MW1 | 3/14/2008 | MW-1 | 1.00 U | 0.320 | 0.200 U | 0.200 U | 0.300 | 0.200 U | 1.00 U | 0.200 U | 2.00 U | 0.200 U | 1.00 U | 1.00 U | 0.200 U | 0.200 U |
| MW2 | 3/14/2008 | MW-2 | 1.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 2.00 U | 0.200 U | 1.00 U | 1.00 U | 0.200 U | 0.200 U |
| MW3 | 3/14/2008 | MW-3 | 1.00 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 2.00 U | 0.200 U | 1.00 U | 1.00 U | 0.200 U | 0.200 U |

BOLD = detection

All results have not been not validated µg/L = micrograms per liter U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

DUP = Field duplicate

Table 9
Volatile Organic Compound Groundwater Analytical Results
SPU- North Recycling and Disposal Station
Seattle, Washington

| | | | p-IsopropyItoluene | sec-Butylbenzene | Styrene | tert-Butylbenzene | Tetrachloroethene | Toluene | Trichloroethene | Trichlorofluoromethane | Vinyl Acetate | Vinyl Chloride | m,p-Xylene | o-Xylene | Total Xylenes |
|---------------|-----------------|-----------|--------------------|------------------|---------|-------------------|-------------------|---------|-----------------|------------------------|---------------|----------------|------------|----------|---------------|
| 14704 14 14 | 0.11 | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method A | | | | | 400 | | 5 | 1,000 | 5 | | | 0.2 | | | 1000 |
| | AR- Federal MCL | | | | 100 | | 5 | 1,000 | 5 | | | 2 | | | 10 |
| | CULs, Carcinoge | | | | 1.5 | | 0.081 | | 0.11 | | | 0.029 | | | |
| | CULs, Noncarcin | 0 | | | 1,600 | | 80 | 640 | 2.4 | 2,400 | 8,000 | 24 | 16,000 | 16,000 | 1,600 |
| Location ID | Sample Date | Sample ID | | | | | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 0.750 | 0.290 | 1.04 |
| MW1 | 3/14/2008 | MW-1 | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 0.750 | 0.300 | 1.05 |
| MW2 | 3/14/2008 | MW-2 | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 0.400 U | 0.200 U | 0.400 |
| MW3 | 3/14/2008 | MW-3 | 0.200 U | 0.200 U | 0.200 U | 0.200 U | 0.810 | 0.200 U | 0.200 U | 0.200 U | 1.00 U | 0.200 U | 0.400 U | 0.200 U | 0.400 |

BOLD = detection

All results have not been not validated

μg/L = micrograms per liter

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) and Federal Primary Maximum Contaminant Level (MCL) from Washington State Department of Ecology, Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

| | | | (3+4)-Methylphenol (m,p-Cresol) | 1,2,4-Trichlorobenzene | 1,2-Dichlorobenzene | 1,2-Dinitrobenzene | 1,2-Diphenylhydrazine | 1,3-Dichlorobenzene | 1,3-Dinitrobenzene | 1,4-Dichlorobenzene | 1,4-Dinitrobenzene | 1-Methylnaphthalene | 2,3,4,6-Tetrachlorophenol |
|--------------------------|-------------|-----------|---------------------------------|------------------------|---------------------|--------------------|-----------------------|---------------------|--------------------|---------------------|--------------------|---------------------|---------------------------|
| | | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method | | | | | | | | | | | | | |
| Groundwater A | | | | 70 | 600 | | | | | | 75 | | |
| MTCA Method | | | 400 | | 700 | | 0.11 | | | 1.4 | | | 400 |
| MTCA Method Location ID | | | 400 | 80 | 720 | | | | | | | 2.4 | 480 |
| | Sample Date | Sample ID | | | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 0.110 U | 1.10 U |
| MW1 | 3/14/2008 | MW-1 | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 0.100 U | 1.00 U |
| MW2 | 3/14/2008 | MW-2 | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.098 U | 0.980 U |
| MW3 | 3/14/2008 | MW-3 | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 0.110 U | 1.10 U |

Notes:

BOLD = detection

All results have not been not validated $\mu g/L = micrograms$ per liter

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) and Federal Primary Maximum Contaminant Level (MCL) from Washington State Department of Ecology, Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

| | | | 2,3,5,6-Tetrachlorophenol | 2,3-Dichloroaniline | 2,4,5-Trichlorophenol | 2,4,6-Trichlorophenol | 2,4-Dichlorophenol | 2,4-Dimethylphenol | 2,4-Dinitrophenol | 2,4-Dinitrotoluene | 2,6-Dinitrotoluene | 2-Chloronaphthalene | 2-Chlorophenol |
|-----------------|-----------------------------------|-----------|---------------------------|---------------------|-----------------------|-----------------------|--------------------|--------------------|-------------------|--------------------|--------------------|---------------------|----------------|
| NATOA NA 41 - 1 | A 01 II | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method | | ICI | | | | | | | | | | | |
| | RAR- Federal M | | | | | 4 | | | | | | | |
| | B CULs, Carcino B CULs, Noncar | | | | 800 | | 24 | 160 | 0.32 | 32 | 16 | | 40 |
| Location ID | Sample Date | Sample ID | | | 000 | | 27 | 100 | 0.02 | 52 | 10 | | |
| MW1 | 3/14/2008 | DUP-1 | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 5.40 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U |
| MW1 | 3/14/2008 | MW-1 | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 5.20 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U |
| MW2 | 3/14/2008 | MW-2 | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 4.90 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U |
| MW3 | 3/14/2008 | MW-3 | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 5.40 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U |

Notes:

BOLD = detection

All results have not been not validated

μg/L = micrograms per liter

--- = no data/not researched

DUP = Field duplicate

U = The analyte is not detected at or above the listed reporting detection limit

| | | | 2-Methylnaphthalene | 2-Methylphenol (o-Cresol) | 2-Nitroaniline | 2-Nitrophenol | 3,3'-Dichlorobenzidine | 3-Nitroaniline | 4,6-Dinitro-2-methylphenol | 4-Bromophenyl-phenylether | 4-Chloro-3-methylphenol | 4-Chloroaniline | 4-Chlorophenyl-phenylether |
|-------------|-----------------------------------|-----------|---------------------|---------------------------|----------------|---------------|------------------------|----------------|----------------------------|---------------------------|-------------------------|-----------------|----------------------------|
| NATOA NA d | A 01 II | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method | | 101 | | | | | | | | | | | |
| | RAR- Federal M | | | | | | 0.10 | | | | | | |
| | B CULs, Carcino B CULs, Noncar | | 32 | 400 | | | 0.19 | | | | | | |
| Location ID | Sample Date | Sample ID | 32 | 400 | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 0.110 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 5.40 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U |
| MW1 | 3/14/2008 | MW-1 | 0.100 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 5.40 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U |
| MW2 | 3/14/2008 | MW-2 | 0.098 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 4.90 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U |
| MW3 | 3/14/2008 | MW-3 | 0.110 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 5.40 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U |

Notes:

BOLD = detection

All results have not been not validated $\mu g/L = micrograms$ per liter

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) and Federal Primary Maximum Contaminant Level (MCL) from Washington State Department of Ecology, Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

| | | | 4-Nitroaniline | 4-Nitrophenol | Acenaphthene | Acenaphthylene | Aniline | Anthracene | Benzidine | Benzo[g,h,i]perylene | Benzyl alcohol | bis(2-Chloroethoxy)methane | bis(2-Chloroethyl)ether |
|--------------------------|-------------|-----------|----------------|---------------|--------------|----------------|---------|------------|-----------|----------------------|----------------|----------------------------|-------------------------|
| | | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method | | 101 | | | | | | | | | | | |
| Groundwater A | | | | | | | 7.7 | | | | | | |
| MTCA Method | | | | | | | 7.7 | 4.000 | 0.00038 | | 0.400 | | 0.04 |
| MTCA Method Location ID | Sample Date | Sample ID | | | 960 | | | 4,800 | 48 | | 2,400 | | |
| | | | | | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 1.10 U | 1.10 U | 0.110 U | 0.110 U | 1.10 U | 0.110 U | 11.0 U | 0.011 U | 1.10 U | 1.10 U | 1.10 U |
| MW1 | 3/14/2008 | MW-1 | 1.00 U | 1.00 U | 0.100 U | 0.100 U | 1.00 U | 0.100 U | 10.0 U | 0.010 U | 1.00 U | 1.00 U | 1.00 U |
| MW2 | 3/14/2008 | MW-2 | 0.980 U | 0.980 U | 0.098 U | 0.098 U | 0.980 U | 0.098 U | 9.80 U | 0.010 U | 0.980 U | 0.980 U | 0.980 U |
| MW3 | 3/14/2008 | MW-3 | 1.10 U | 1.10 U | 0.110 U | 0.110 U | 1.10 U | 0.110 U | 11.0 U | 0.011 U | 1.10 U | 1.10 U | 1.10 U |

Notes:

BOLD = detection

All results have not been not validated $\mu g/L = micrograms per liter$

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) and Federal Primary Maximum Contaminant Level (MCL) from Washington State Department of Ecology, Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

| | | | bis(2-Chloroisopropyl)ether | bis(2-Ethylhexyl)phthalate | bis-2-Ethylhexyladipate | Butylbenzylphthalate | Carbazole | Dibenzofuran | Diethylphthalate | Dimethylphthalate | Di-n-butylphthalate | Di-n-octylphthalate | Fluoranthene |
|--------------------------|-------------|-----------|-----------------------------|----------------------------|-------------------------|----------------------|-----------|--------------|------------------|-------------------|---------------------|---------------------|--------------|
| | | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method | | 101 | | | | | | | | | | | |
| Groundwater A | | | | 6 | | | 4.4 | | | | | | |
| MTCA Method | | | 220 | 6.3 | | 2 200 | 4.4 | 0.22 | 12.000 | 16.000 | | 220 | 640 |
| MTCA Method Location ID | Sample Date | Sample ID | 320 | 320 | | 3,200 | | 0.32 | 13,000 | 16,000 | | 320 | 640 |
| | | | 4.40.11 | 4.40.11 | 4.40.11 | 4.40.11 | 4.40.11 | 4.40.11 | 4.40.11 | 4.40.11 | | 4.40.11 | 0.440.11 |
| MW1 | 3/14/2008 | DUP-1 | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 2.30 | 1.10 U | 0.110 U |
| MW1 | 3/14/2008 | MW-1 | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 2.20 | 1.00 U | 0.100 U |
| MW2 | 3/14/2008 | MW-2 | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.098 U |
| MW3 | 3/14/2008 | MW-3 | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 0.110 U |

Notes:

BOLD = detection

All results have not been not validated $\mu g/L = micrograms$ per liter

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) and Federal Primary Maximum Contaminant Level (MCL) from Washington State Department of Ecology, Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

| | | | Fluorene | Hexachlorobenzene | Hexachlorobutadiene | Hexachlorocyclopentadiene | Hexachloroethane | Isophorone | Naphthalene | Nitrobenzene | N-Nitrosodimethylamine | N-Nitroso-di-n-propylamine | N-Nitrosodiphenylamine |
|--------------------------|-----------------|-----------|----------|-------------------|---------------------|---------------------------|------------------|------------|-------------|--------------|------------------------|----------------------------|------------------------|
| | | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method | | 101 | | | | | | | | | | | |
| | RAR- Federal M | | | 1 | 0.50 | 50 | | | | | | | |
| | B CULs, Carcino | | | 0.055 | 0.56 | 40 | 3.1 | 0.46 | 400 | | 0.00086 | | |
| MTCA Method Location ID | Sample Date | Sample ID | 640 | 13 | 1.6 | 48 | 8 | 1,600 | 160 | 4 | | | |
| | | | | | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 0.110 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 0.110 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U |
| MW1 | 3/14/2008 | MW-1 | 0.100 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U | 0.100 U | 1.00 U | 1.00 U | 1.00 U | 1.00 U |
| MW2 | 3/14/2008 | MW-2 | 0.098 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U | 0.098 U | 0.980 U | 0.980 U | 0.980 U | 0.980 U |
| MW3 | 3/14/2008 | MW-3 | 0.110 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U | 0.110 U | 1.10 U | 1.10 U | 1.10 U | 1.10 U |

Notes:

BOLD = detection

All results have not been not validated

μg/L = micrograms per liter

U = The analyte is not detected at or

--- = no data/not researched

above the listed reporting detection limit

DUP = Field duplicate

| | | | Pentachlorophenol | Phenanthrene | Phenol | Pyrene | Pyridine | Benzo[a]anthracene | Benzo[a]pyrene | Benzo[b]fluoranthene | Benzo[k]fluoranthene | Chrysene | Dibenz[a,h]anthracene | Indeno[1,2,3-cd]pyrene | Total TEF |
|----------------|-------------|-----------|-------------------|--------------|---------|---------|----------|--------------------|----------------|----------------------|----------------------|----------|-----------------------|------------------------|-----------|
| | | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method | | | | | | | | | 0.1 | | | | | | 0.1 |
| Groundwater Al | | | 1 | | | | | | 0.2 | | | | | | 0.2 |
| MTCA Method I | | | 0.73 | | 4.000 | | | 0.012 | 0.012 | 0.012 | 0.012 | 0.012 | 0.012 | 0.012 | 0.012 |
| MTCA Method | | | 480 | | 4,800 | 480 | 8 | | | | | | | | |
| Location ID | Sample Date | Sample ID | | | | | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 5.40 U | 0.110 U | 1.10 U | 0.110 U | 1.10 U | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.006 |
| MW1 | 3/14/2008 | MW-1 | 5.20 U | 0.100 U | 1.00 U | 0.100 U | 1.00 U | 0.010 U | 0.010 U | 0.010 U | 0.010 U | 0.010 U | 0.010 U | 0.010 U | 0.008 |
| MW2 | 3/14/2008 | MW-2 | 4.90 U | 0.098 U | 0.980 U | 0.170 | 0.980 U | 0.010 U | 0.010 U | 0.010 U | 0.010 U | 0.010 U | 0.010 U | 0.010 U | 0.008 |
| MW3 | 3/14/2008 | MW-3 | 5.40 U | 0.110 U | 1.10 U | 0.110 U | 1.10 U | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.011 U | 0.006 |

Notes:

BOLD = detection

All results have not been not validated TEF = Toxicity Equivalency Factor calculated in accordance with WAC 173-340-708(e)

 μ g/L = micrograms per liter U = The analyte is not detected at or

--- = no data/not researched above the listed reporting detection limit

DUP = Field duplicate

Table 11
Dissolved Metal Groundwater Analytical Results
SPU- North Recycling and Disposal Station
Seattle, Washington

| | | | Arsenic | Barium | Cadmium | Chromium | Lead | Mercury | Selenium | Silver |
|---------------|-----------------|-----------|---------|--------|---------|----------|--------|---------|----------|--------|
| | A Method A CUL | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method | A CUL | | 5 | | 5 | 50 | 15 | 2 | 50 | |
| Groundwater A | RAR- Federal M | CL | 20 | 2,000 | 5 | 100 | 15 | 2 | 50 | |
| MTCA Method | B CULs, Carcino | gen | 0.058 | | | | | | | |
| MTCA Method | B CULs, Noncard | cinogen | 4.8 | 3,200 | 8 | | | 4.8 | 80 | 80 |
| Location ID | Sample Date | Sample ID | · | · | · | · | | · | | |
| MW1 | 3/14/2008 | DUP-1 | 3.00 U | 31.0 | 4.00 U | 10.0 U | 1.00 U | 0.500 U | 5.00 U | 10.0 U |
| MW1 | 3/14/2008 | MW-1 | 3.00 U | 33.0 | 4.00 U | 10.0 U | 1.00 U | 0.500 U | 5.00 U | 10.0 U |
| MW2 | 3/14/2008 | MW-2 | 3.00 U | 25.0 U | 4.00 U | 10.0 U | 1.00 U | 0.500 U | 5.00 U | 10.0 U |
| MW3 | 3/14/2008 | MW-3 | 3.00 U | 36.0 | 4.00 U | 10.0 U | 1.00 U | 0.500 U | 5.00 U | 10.0 U |

BOLD = detection

All results have not been not validated µg/L = micrograms per liter

--- = no data/not researched

DUP = Field duplicate

MTCA Method A and B Cleanup Levels (CULs) and Federal Primary Maximum Contaminant Level (MCL) from Washington State Department of Ecology, Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

U = The analyte is not detected at or

Table 12 Pesticide Groundwater Analytical Results SPU- North Recycling and Disposal Station Seattle, Washington

| | | | 4,4'-DDD | 4,4'-DDE | 4,4'-DDT | Aldrin | alpha-BHC | alpha-Chlordane | beta-BHC | delta-BHC | Dieldrin | Endosulfan I | Endosulfan II | Endrin | Endrin Aldehyde | Endrin Ketone | Endsulfan Sulfate | gamma-BHC | gamma-Chlordane | Heptachlor | Heptachlor Epoxide | Methoxychlor | Toxaphene |
|-------------|-----------------|-----------|----------|----------|----------|---------|-----------|-----------------|----------|-----------|----------|--------------|---------------|---------|-----------------|---------------|-------------------|-----------|-----------------|------------|--------------------|--------------|-----------|
| | | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Metho | d A CUL | | | | 0.3 | | | | | | | | | | | | | | | | | | |
| Groundwater | ARAR- Federal | MCL | | | | | | | | | | | | 2 | | | | | | 0.4 | 0.2 | 40 | |
| MTCA Metho | d B CULs, Carci | nogen | 0.36 | 0.26 | 0.26 | 0.0026 | | | | | 0.0055 | | | | | | | | | 0.019 | 0.0048 | | 0.08 |
| MTCA Metho | d B CULs, Nonc | arcinogen | | | 8 | 0.24 | | | | | 0.8 | 96 | | 4.8 | | | | | | 8 | 0.1 | 80 | |
| Location ID | Sample Date | Sample ID | | | | | | | | | | | | | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.019 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.010 U | 0.048 U |
| MW1 | 3/14/2008 | MW-1 | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.023 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.006 U | 0.011 U | 0.056 U |
| MW2 | 3/14/2008 | MW-2 | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.020 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.010 U | 0.050 U |
| MW3 | 3/14/2008 | MW-3 | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.020 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.005 U | 0.010 U | 0.051 U |

Notes:

BOLD = detection

All results have not been not validated $\mu g/L = micrograms per liter$

U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

DUP = Field duplicate

Table 13
Herbicide Groundwater Analytical Results
SPU- North Recycling and Disposal Station
Seattle, Washington

| | | | 2,4,5-T | 2,4,5-TP (Silvex) | 2,4-D | 2,4-DB | Dalapon | Dicamba | Dichlorprop | Dinoseb | MCPA | MCPP | Pentachlorophenol |
|-----------------|--------------------|-----------|---------|-------------------|---------|---------|---------|---------|-------------|---------|--------|--------|-------------------|
| | | | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) | (µg/L) |
| MTCA Method A | CUL | | | | | | | | | | | | |
| Groundwater ARA | AR- Federal MCL | | | | | | | | | | | | |
| MTCA Method B | CULs, Carcinogen | | | | | | | | | | | | 0.73 |
| MTCA Method B | CULs, Noncarcinoge | n | | | 160 | 130 | 240 | 480 | | | 16 | | 480 |
| Location ID | Sample Date | Sample ID | | | | | | | | | | | |
| MW1 | 3/14/2008 | DUP-1 | 0.024 U | 0.024 U | 0.470 U | 0.024 U | 0.230 U | 0.024 U | 0.024 U | 0.024 U | 4.70 U | 4.70 U | 0.010 U |
| MW1 | 3/14/2008 | MW-1 | 0.023 U | 0.023 U | 0.460 U | 0.023 U | 0.230 U | 0.023 U | 0.023 U | 0.023 U | 4.60 U | 4.60 U | 0.009 U |
| MW2 | 3/14/2008 | MW-2 | 0.023 U | 0.023 U | 0.460 U | 0.023 U | 0.220 U | 0.023 U | 0.023 U | 0.023 U | 4.60 U | 4.60 U | 0.009 U |
| MW3 | 3/14/2008 | MW-3 | 0.023 U | 0.023 U | 0.450 U | 0.023 U | 0.220 U | 0.023 U | 0.023 U | 0.023 U | 4.50 U | 4.50 U | 0.009 U |

BOLD = detection

All results have not been not validated $\mu g/L = micrograms per liter$

U = The analyte is not detected at or above the listed reporting detection limit

--- = no data/not researched

DUP = Field duplicate

Table 14 Polychlorinated Biphenyl Groundwater Analytical Results SPU- North Recycling and Disposal Station Seattle, Washington

| | | | Aroclor 1016 | Aroclor 1221 | Aroclor 1232 | Aroclor 1242 | Aroclor 1248 | Aroclor 1254 | Aroclor 1260 | Aroclor 1262 | Aroclor 1268 | Total Aroclors |
|-----------------|------------------|-----------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|--------------|----------------|
| | | | (µg/L) |
| MTCA Method A | CUL | | | - | | | | | | | | 0.10 |
| Groundwater ARA | AR- Federal MCL | | | | | | | | | | | 0.50 |
| MTCA Method B | CULs, Carcinogen | | | | | | | | | | | 0.044 |
| MTCA Method B | CULs, Noncarcino | gen | | - | | | | | | | | |
| Location ID | Sample Date | Sample ID | | | | · | | | · | | | |
| MW1 | 3/14/2008 | DUP-1 | 0.058 U | ND |
| MW1 | 3/14/2008 | MW-1 | 0.050 U | ND |
| MW2 | 3/14/2008 | MW-2 | 0.048 U | ND |
| MW3 | 3/14/2008 | MW-3 | 0.048 U | ND |

Notes:

BOLD = detection

ND = Not detected above laboratory detection limits

U = The analyte is not detected at or above the listed reporting detection limit

All results have not been not validated

 μ g/L = micrograms per liter

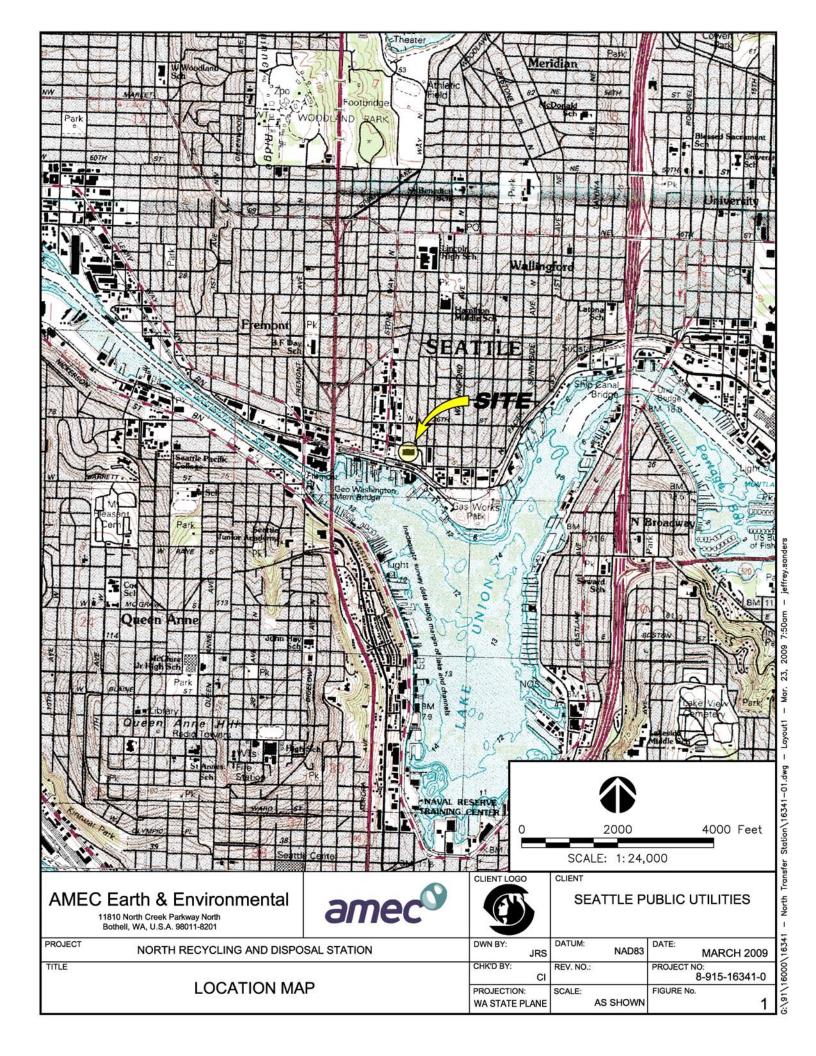
--- = no data/not researched

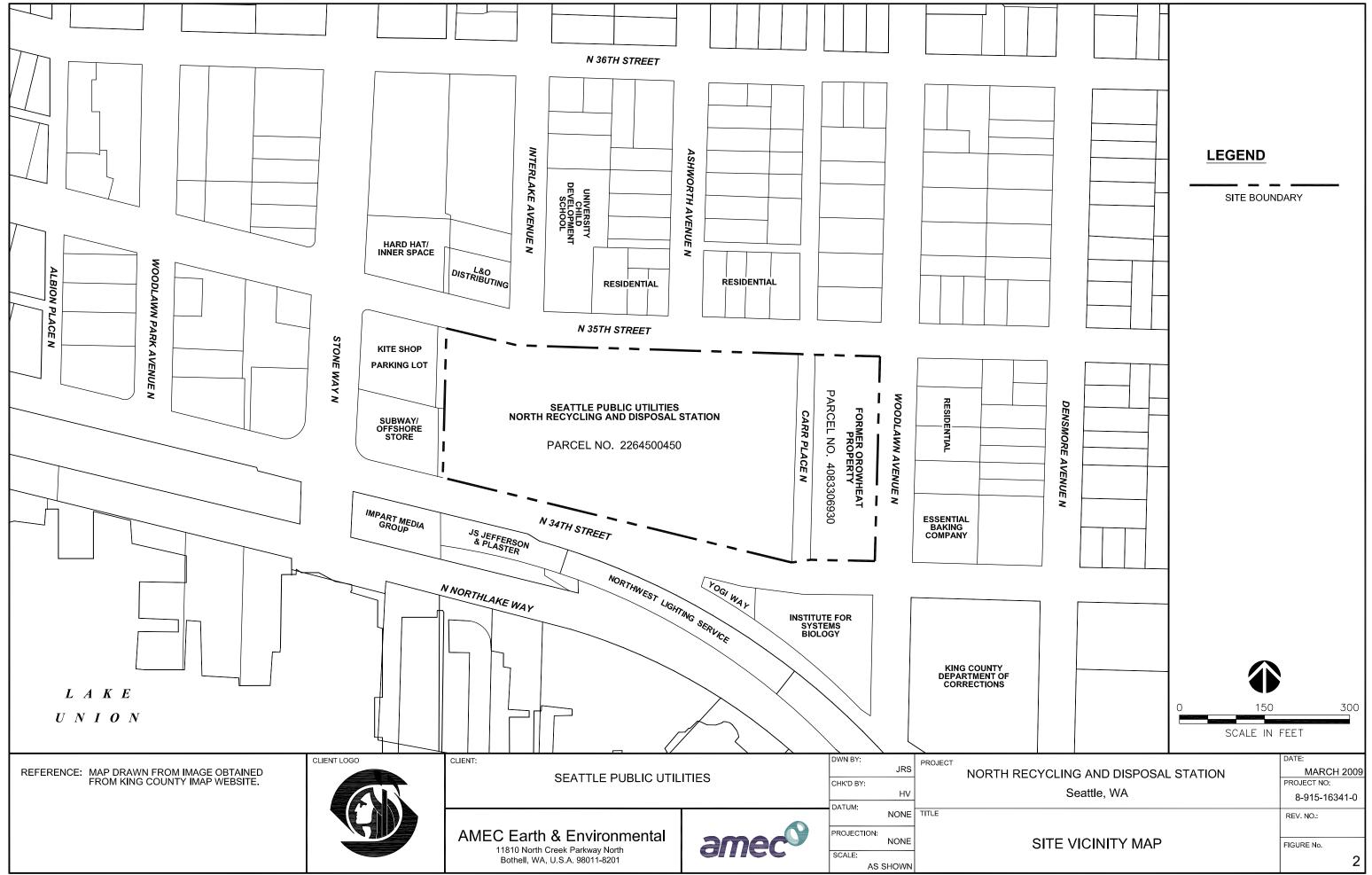
DUP = Field duplicate

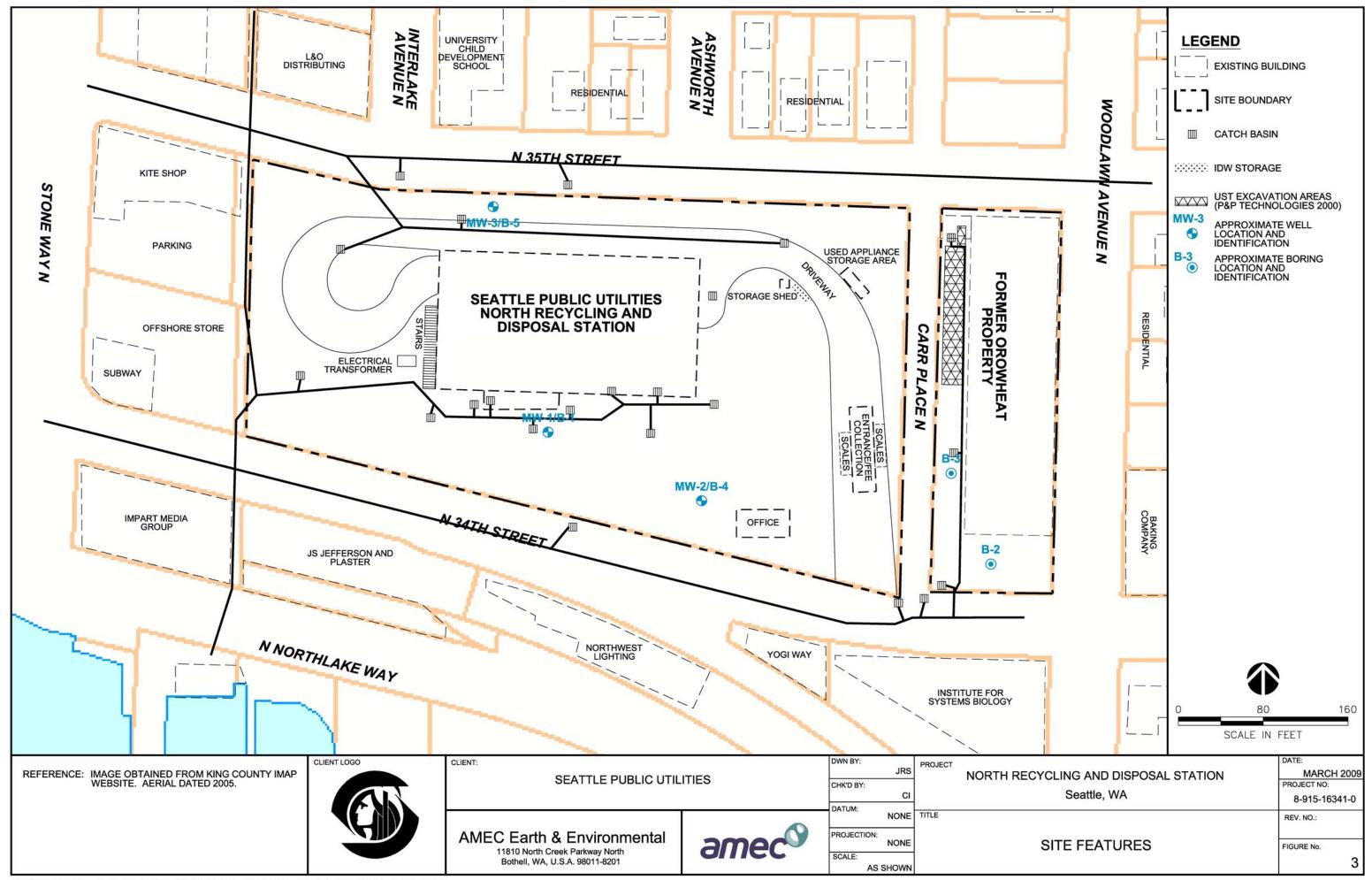
MTCA Method A and B Cleanup Levels (CULs) and Federal Primary Maximum Contaminant Level (MCL) from Washington State Department of Ecology, Cleanup Levels and Risk Calculations (CLARC) database (April 17, 2008)

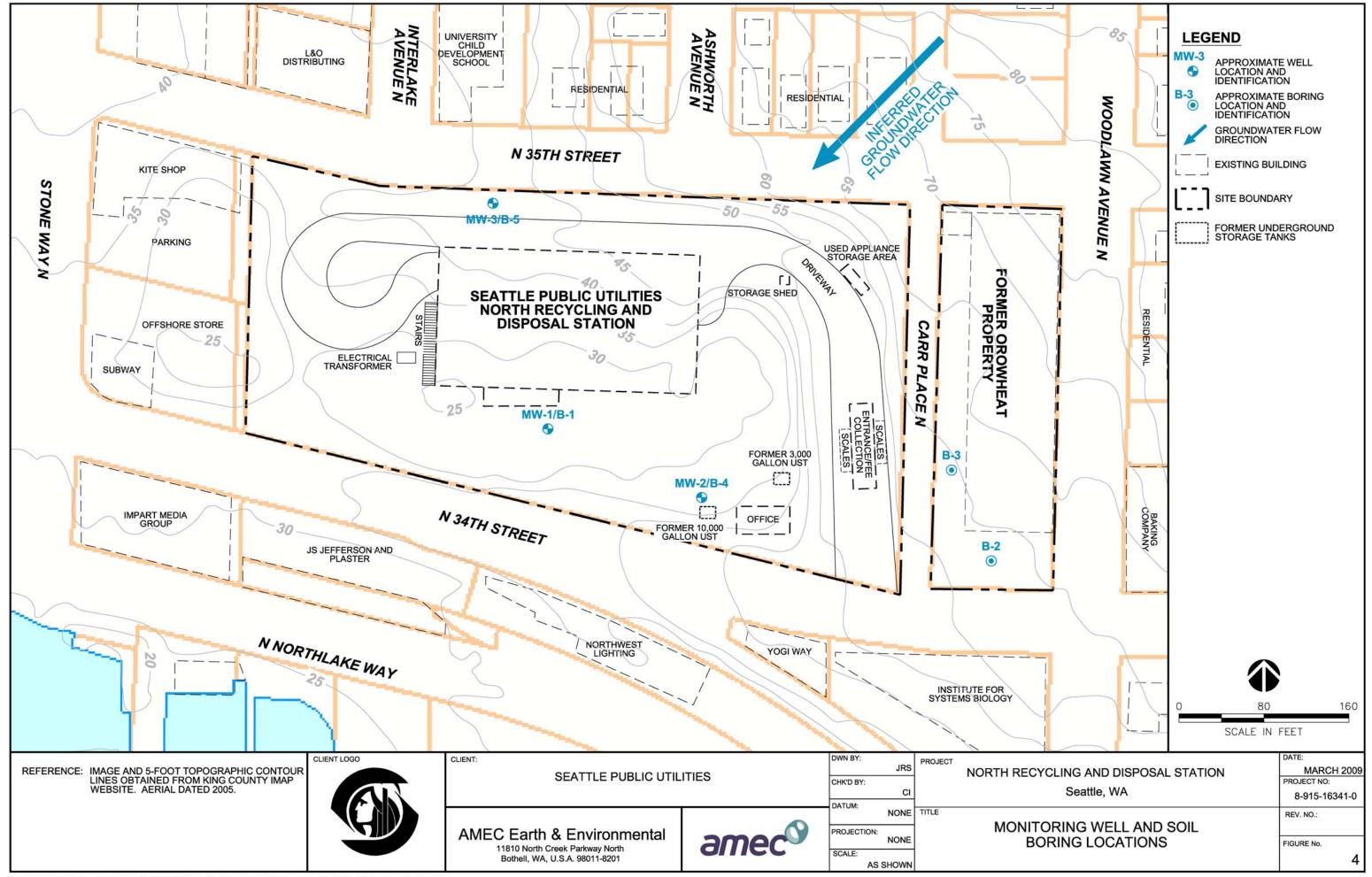


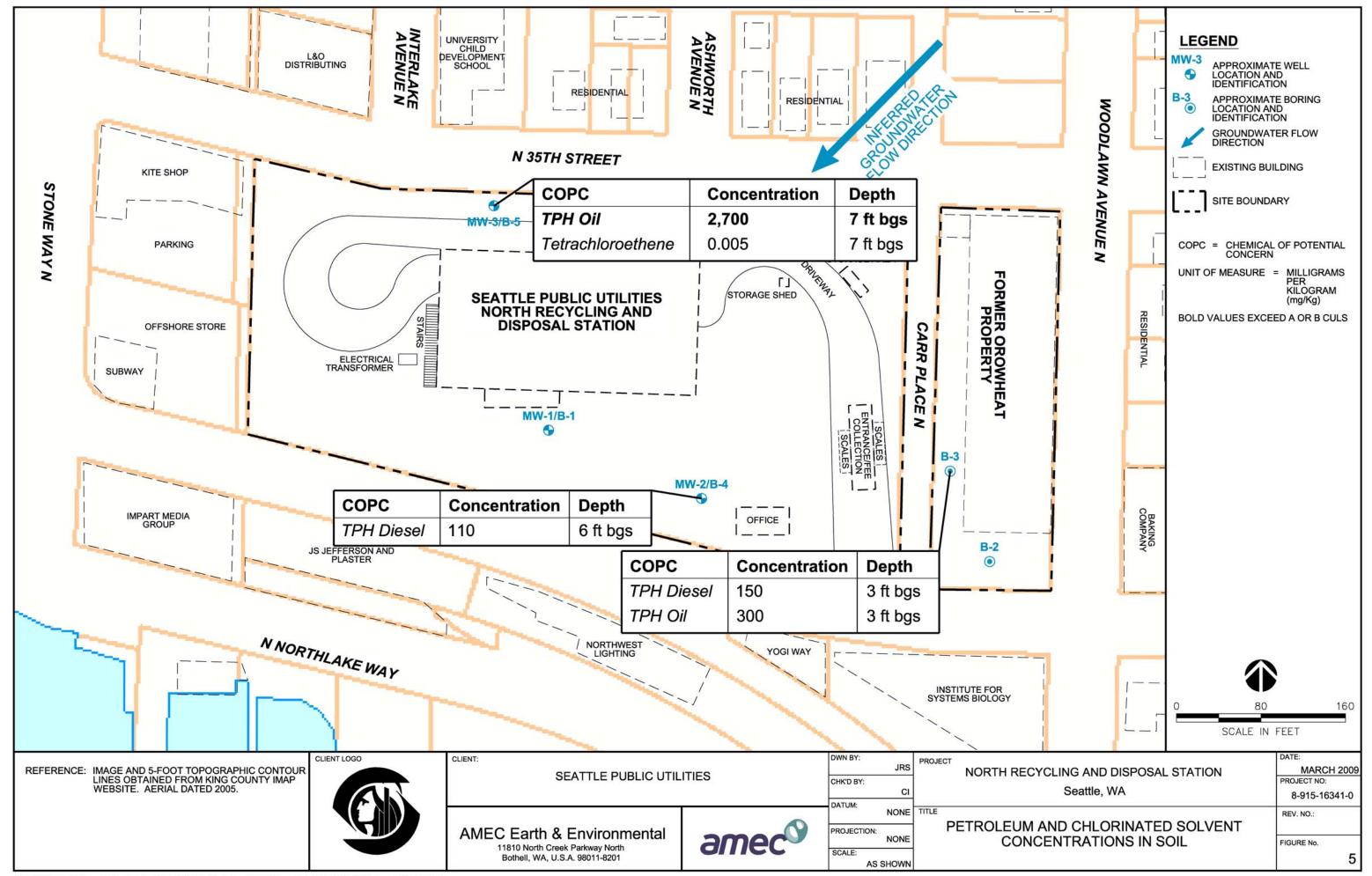
FIGURES

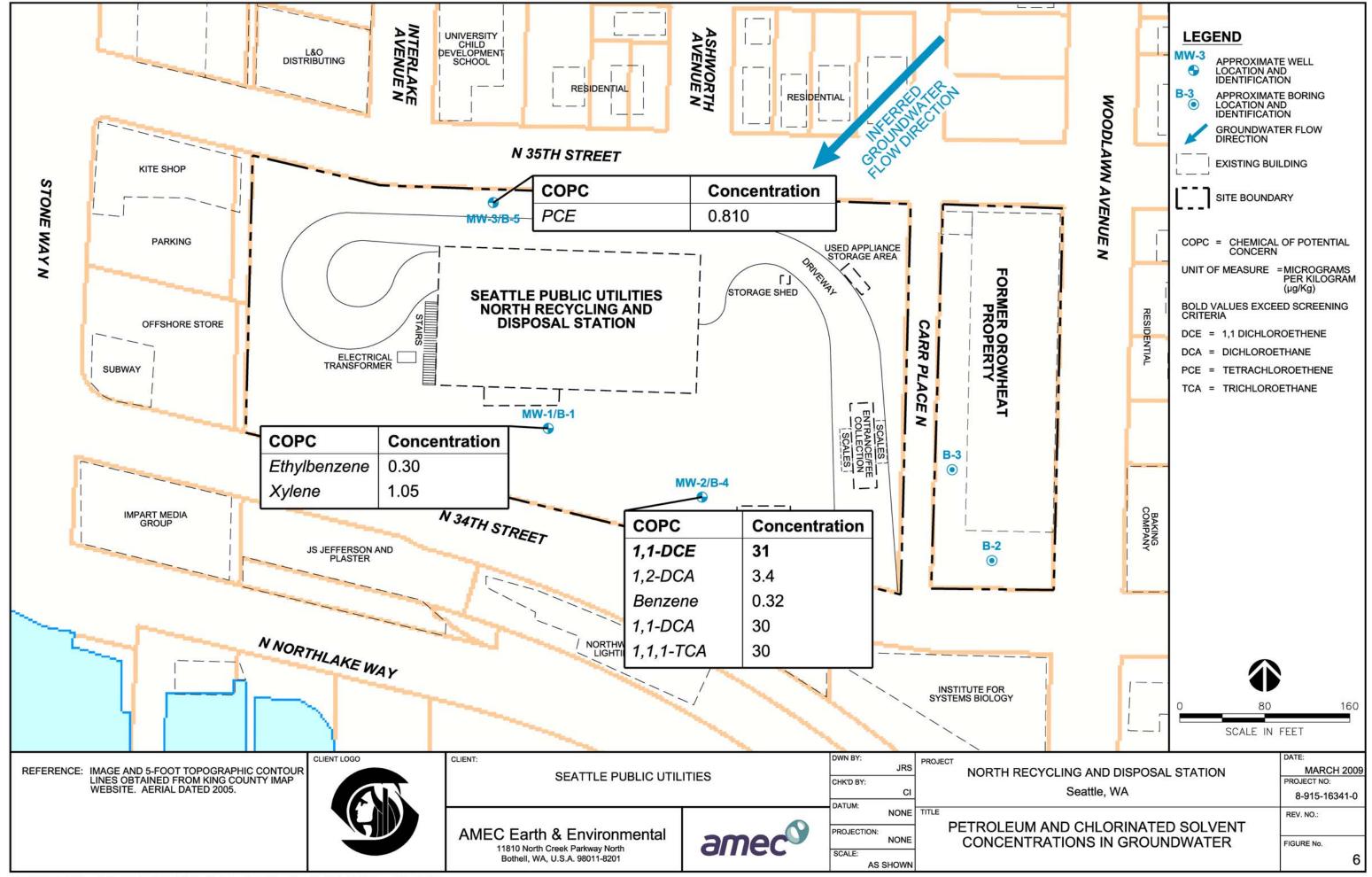














APPENDIX A

Monitoring Well Development Logs

| | | _ | PROJECT | | | | | ELL NO. | | | | | |
|---|--------------------------------|---|-----------------|------------------------|--|--|--------------|---------|--|---|-----------------|---------------------|------------|
| , | WELL D | EVELC | PME | NT LO | G | JOB NO. | | SITE | : | P | REPARED 8 | | |
| METHO | D VERPUMPAGE | | | NITIAL WAT | ER I EVEL | 7.(FC | • , | R | EMARKS: | n Obse | **** | | |
| | AILER | | | INAL WATE | | | | | P20 = | | | | |
| | URGE LOCK | | | CAPACITY (GALLONS/I | OF CASING | | | | IE BETWEE (GALLONS ASSUMING | /LINEAR F | | | |
| Al | R LIFT | | | _ | = 0.16 = 0.65 | 5 2" CASING AND 6" HOLE - 0.52 | | | | | | | |
| | THER | | | 6* = | = 1.47 | 2" CASING AND 8" HOLE - 0.98 4" CASING AND 10" HOLE = 1.37 4" CASING AND 12" HOLE - 2.09 | | | | | | | |
| HOLE DIAMETER d _n = d _w WELL CASING INSIDE DIAMETER OUTSIDE DIAMETER d _w OD = booder DEPTH TO: WATER LEVEL H DEPTH TO: WATER LEVEL DEPTH TO: WATER LEVEL | | | | CAS | WELL VOLUME CALCULATION: CASING VOLUME = $V_c = \Pi \left(\frac{d_b ID}{2} \right)^2 (TD - H) = 3.14 \left(\frac{d_b OD}{2} \right)^2 \left(\frac$ | | | | | | | | |
| BASE (| OF SEAL OF WELL TER PACK | H = <u>7.1</u> S = TD = <u>14.4</u> P = | _ | screen Interval | o ! | | | | $(*if S > $ $= 3.14 \left[\left({2} \right) \right]$ | H, use S: $- \int_{0}^{2} -\left(\frac{1}{2}\right)^{2} dt$ | if S < H, use 1 | | _ |
| | DE | VELOPMENT LO | G: | | CUMULAT | TVE | | | ER QUALITY | | | | 7 |
| DATE | TIME BEGIN/END | METHOD | ELAPSED TIME | FLOW RATE (gpm) | REMOVE GALLON | :D 0: | TEMP | CONDUC | | REDOX | TURBID- ITY | COMMENTS | |
| 3.(2. a r | 1313 | HISCH PAN | | | | 2.10 | (1.41 | 0.49 | 1 13.33 | 281 | 299+ | Silty | |
| a | 1715 | a | Z min | | 4 72 | દૈ. વિબ્ર⊊ | 11.84 | 0.665 | 10.32 | 241 | 999+ | silly pamped day | 17 min ADR |
| | 1344 | 11 | 3 mm | | | 2.35 | 12.34 | 0.653 | 10.17 | 237 | 357 TA | Silty Pumper any | EFT Acuse |
| | 1430 | t, | 3 min | | APROY. | 7.30 | 12-13 | 0.721 | 10.54 | 218 | 911 + | pumped Day | <u>/</u> |
| | | | | | | | | | | | | | _ |
| | | | | | | | | | <u> </u> | <u> </u> | | | _ |
| | | | | | | | | | | | | | _ |
| | | | | | | | | | | | | | _ |
| | <u> </u> | | | | | | | | | | | | - |
| | | | | | | | | | | | | | _ |
| | | | | | | | | | | | | | |
| 2 | | | | | | | | | | | | | |
| | | | | | | | | | | | | | |
| | | | | | | | | | _ | | | | 1 |
| | | | | | | | | | | | | | - |

* = Dissolved Oxygen

| | WELL DEVELOPMENT LOG | | | | _ | PROJECT | | | WI | WELL NO. | | | | |
|--|----------------------|---|-----------------|-----------------------|--------------------------------------|---|--------|----------|---------------------------------------|---|----------------------------------|---|---------------------|-------------|
| 1 | WELL D | EVELO | PMEN | IT LO | G | JOE | B NO. | · | SITE | | PF | REPARED E | | 1 |
| METHO | D VERPUMPAGE | | IN | ITIAL WATE | ER LEVEL_ | 5. | 44 | | | MARKS: | ж. | | | = |
| B/ | AILER | | FI | NAL WATE | R LEVEL | | | | 1 | 720 = 0 | may o. | • | | |
| | JRGE LOCK | | I | CAPACITY SALLONS/L | | _ | | | (0 | ALLONS | N CASING LINEAR F 40% PORC | AND HOLI OOT) | E | |
| AI | R LIFT | | | _ | 0.16 | | | | | | ND 6" HOLI | | | |
| 0- | THER | | | | 2" CASING AND 8" HOLE - 0.98 1.47 | | | | | | | | | |
| HOLE DI | AMETER | d _h = | _ | ٠ | WELL V | VOLU | ME CA | LCULAT | | | | | | 1 |
| WELL CA INSIDE OUTSIE DIAMETE | DIAMETER DE | $d_{w}ID = 2$ $d_{w}OD = \frac{2}{100}$ $d_{w}OD = \frac{2}{100}$ $CASING VOLUME = V_{c} = \Pi \left(\frac{d_{w}ID}{2}\right)^{2} (TD - H) = 3.14 \left(\frac{1}{2}\right)^{2} \left($ | | | | | | = | | | | | | |
| DEPTH T WATER | | H = <u>5.4</u> | | | FILT | ER PA | ACK PO | ORE VOLU | ME = Vı= | $\Pi\left[\left(\frac{\mathrm{dh}}{2}\right)^2-\left(\frac{\mathrm{dh}}{2}\right)^2\right]$ | $\left(\frac{d_wOD}{2}\right)^2$ | TD – (S or H | *)(P) = | |
| | | s = | _ | SCREENED INTERVAL | . | | | | | • | | fS <h, td="" use<=""><td>•</td><td></td></h,> | • | |
| EST. FIL | TER PACK | TD = <u>(7.4</u> | | d, - | | | | | = | 3.14 | -)2-(| -)²] |)() = | |
| POROSIT | ΙΥ | P = | - | | тот. | AL W | ELL V | OLUME = | VT = Vc + \ | /f = | =_ | ft.³ x 7 | 7.48 =gal. | |
| | DE | VELOPMENT LO | G: | | CUMULAT | | | • | WATER | QUALITY | | | | 1 |
| DATE | TIME BEGIN/END | METHOD | ELAPSED TIME | FLOW RATE (gpm) | REMOVE GALLON | :D | рΗ | TEMP | CONDUC- TIVITY | D.O.* | REDOX | TURBID- ITY | COMMENTS | |
| 3.12.0F | 1509 | HISCH FROM PUMP | Zmin | | APPOF 4 GM | <u>, </u> | 7.03 | 11.41 | 0.463 | 11.00 | 209 | 1177 | Silty PUMPED DAY | 3 m R€-0 |
| k | 1540 | 11 | Zmin | | APROX 4 GAL. | | 7.14 | 12.92 | 0.571 | 10.50 | Ito | 454 + | Silly Funces Day | 1.5 |
| n | 1420 | 2, | 2mh | | Annux. 4 Gal | | 7.10 | 12.43 | 0.560 | 10.43 | 180 | 500 | pumper Day | |
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* = Dissolved Oxygen

| _ | WELL DEVELOPMENT LOG | | | | | PROJECT | | | | | WI | WELL NO. | | | 1 |
|---|----------------------|--------------------------------|--|-----------------------|--|--|-------|-----------------------|---|------------|-------------------------------------|---|---------------------|-----|----------|
| 1 | WELL D | EVELO | PME | NT LO | G | JOE | B NO. | | SITE | | | REPARED B T. PERAZU | Y | |] |
| METHO O | D VERPUMPAGE | | ı | NITIAL WAT | ER LEVEL_ | 13. | 27 | | | MARKS: | n),0 ppm. | | | | |
| | AILER JRGE | | <u>_ </u> | INAL WATE | R LEVEL | G | | | | , | | AND HOLE | : | | - |
| Bl | OCK | | | (GALLONS/L | INEAR FOO | | | | | | | | | | |
| Al | R LIFT | | | _ | = 0.16 = 0.65 | 2" CASING AND 6" HOLE - 0.52 2" CASING AND 8" HOLE - 0.98 | | | | | | | | | |
| oʻ | THER | | | 6" = | 4" CASING AND 10" HOLE = 1,37 4" CASING AND 12" HOLE - 2.09 | | | | | | | | | | |
| HOLE DI | | d _h = | _ | GROUND SURFACE | WELL | VOLU | ME CA | LCULATI | ON: | | | | | | |
| WELL CA | ASING DIAMETER | $d_{\text{MID}} = \frac{2}{2}$ | - 1 | | · | | | | 4.ID},' | > | $\int_{0}^{\infty} \int_{0}^{2} dt$ | , | | | |
| OUTSIC DIAMETE | | d = | _ | # * s | CAS | ING V | /OLUM | IE = V _c = | $\Pi\left(\frac{\mathrm{d-ID}}{2}\right)^2$ (TI | O-H) = 3.1 | 4(<u>2</u>)(_ |] = | · | | |
| DEPTH TO: WATER LEVEL $H = \underbrace{13.27} \text{FILTER PACK PORE VOLUME} = V_f = \Pi \left[\left(\frac{d_h}{2} \right)^2 - \left(\frac{d_w OD}{2} \right)^2 \right] \text{TD} - (S \text{ or } H^*)(P) = \underline{\hspace{1cm}}$ | | | | | | _ | | | | | | | | | |
| BASE OF SEAL S = SCREENED | | | | | | | | | | | | fS <h, td="" use<=""><td></td><td></td><td></td></h,> | | | |
| BASE OF WELL EST. FILTER PACK POROSITY P = | | | | | | | | = | 3,14 | | -)'} |)() = . | | | |
| P = | | | | | тот | AL W | ELL V | OLUME = | $VT = V_C + V$ | /f = | + <u></u> =_ | ft.³ x 7 | x 7.48 = gal. | | |
| DEVELOPMENT LOG: | | | | CUMULAT WATER | | | | WATE | QUALITY | | | COMMENTS | | | |
| DATE | TIME BEGIN/END | METHOD | ELAPSED TIME | FLOW RATE (gpm) | REMOVE GALLON | D | рН | TEMP | CONDUC- TIVITY | D.O.* | REDOX | TURBID- ITY | | | |
| 61.12-6> | 1700 | HZLH FACE (RAPET | 2 | | Annov | <u>'</u> | 5.74 | 12.34 | 0.528 | 11.89 | 242 | 919+ | Silly to CI | | /0 21 |
| 4 | 1730 | ч | 2 min | | 4 GAR | | 5.81 | 12.01 | 0.547 | 11.08 | 246 | 9991 | L SILLS PUMPED ! | pηγ | r z |
| l, | 1755 | er . | 2-4 | | APROT 4 CM | | 5.jo | 12.00 | 0.571 | 11.02 | 245 | 350 | elean pumpers | ومه | |
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* = Dissolved Oxygen



APPENDIX B

Groundwater Sample Collection Logs

pg 1 of 2

| amec [©] | LOW-FI | LOW GF | ROUND | WATE | ER SAMP | LING LO | G | | |
|---|--|---|--|--|---|--|---|---|-----------------|
| WELL NO: MUD-1 DATE 3. 14. CO OVAPID READING WHEN WELL STATIC WATER LEVEL (TOC): | LOCATION: S D TIME: L L OPENED: | 14. んり | 001 00 4 92 | CLIMATI DEPTH 1 | C CONDITIONS: TO PRODUCT (TO DEPTH OF WELL | DC): | NA 14,2 | 1 Hoxyric 112193 | HIO |
| METHOD OF I | REMOVAL: 💪 | w for | ושקכ | HUKE | CPURPURP | ING RATE: | <u>Уз д</u> , | <u>lm</u> in | |
| WELL DTW PURGE 7,55 DATA 8,42 7,48 7,55 7,61 7,60 7,73 7,84 7,91 7,95 7,99 8,02 | 1423 1423 1423 1425 1427 1431 1433 1435 1437 1437 1441 1443 | Removed Removed 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 | 11.35 11.35 11.35 11.41 11.35 11.19 11.29 11.20 11.00 11.00 11.00 11.00 | 661 683 698 101 688 100 705 708 709 696 7.13 | Sp. Cond. (mS/cm) 0.777 0.181 0.194 0.795 0.798 0.795 0.838 0.831 0.843 0.847 0.847 | Turbidity (NTU) ent 199 ent 198 749 117 97.0 88.7 207 84.7 | 10.33 10.33 10.29 12.23 10.70 11.12 11.13 11.44 9.68 8.48 10.33 9.07 | 134 134 133 139 142 143 148 144 145 140 137 138 134 | DO du NOT C |
| SAMPLE WITHDRAWAL METHO SAMPLE NUMBER(S) AND TIME | | wer | 4 DL | 1P-T | | | BY: 115 | 01 | <u> </u> |
| NOTES: DD LOTE LAB ANALYSIS PARAMETERS HCI HCI HCI HCI HCI HCI HCI HC | | 10MLY atives: LYVV | DWT TWCA HILLO | PCS OH-(a B/M exed | ax (HC) | Quined s) , Du esticulo 15 (non | StPH-S | How 14 Dx4 O(| 13.746 11(1) |
| NUMBER AND TYPES OF SAMP | PLE CONTAINE -500 M | RS USED: | 10- | 40 c 8 | nluiz -1000 | 115, 4 1001-9 | -500 M | 1100 | <u>55</u> |
| DECON. PROCEDURES: SC OHNC EQU | TOPO C | H Q | OF: | | H20 | erla | neter, | 711 | <u> </u> |
| SAMPLES DELIVERED TO: | 1514 1314 | .08 | | TRANSF | PORTER: | 746 | | | _ |

CAPACITY OF CASING (GALLONS/JINEAR FOOT) 2' - 0.16 + 4' - 0.65 + 6' - 1 47 + 6' - 2 61 + 10' - 4.08 + 12' - 5 57

Pumpon 14B

Dump off 1449

continuation of MW-1

pglofZ

| amec LOW-FLOW GROUNDWATER SAMPLING LOG | | | | | | | | | | | |
|--|-------------------|----------------------|----------------------------|-------------------------|--------------------|-------------------------------|---------------------------|----------------------------------|-------------|--|--|
| WELL NO: | Mial | LOCATION: | | | | | PROJE | CT NO: | ,-, | | |
| DATE: | | TIME: | | | CLIMATION | CONDITIONS: | | | | | |
| OVA/PID RE | ADING WHEN WEL | L OPENED: | | DEPTH TO PRODUCT (TOC): | | | | | | | |
| STATIC WAT | TER LEVEL (TOC): | | | | TOTAL D | EPTH OF WELL (| TOC): | | | | |
| | METHOD OF | REMOVAL: | | | | PUMPIN | G RATE: | | | | |
| WELL PURGE DATA | 8:06 8:10 | Time 1445 1447 | Gallons Removed 4/49 | Temp. (C°) (C°) (11.13 | рн 7,14 7,15 | Sp. Cond. (mS/cm) . 850 | Turbidity (NTU) S.D. 77.8 | 00 4 (mgA) 4 9.93 10.29 | (mv) 136 | | |
| | | | | | | | | | | | |
| | THDRAWAL METHO | | | | | | SAMPLED | BY: | | | |
| SAMPLE NU | IMBER(S) AND TIMI | ±: | | | | | | | | | |
| NOTES: | | | | | · | | ••• | | | | |
| LAB ANALYS | SIS PARAMETERS | AND PRESERV | ATIVES: | - | . | | | <u></u> | | | |
| NUMBER AN | ND TYPES OF SAM | PLE CONTAINE | RS USED: | | | | | | | | |
| DECON. PR | OCEDURES: | | | | | - | | | | | |
| SAMPLES D | ELIVERED TO: | | | | TRANSP | ORTER: | | | | | |

CAPACITY OF CASING [GALLOHS/LINEAR FOOT] 2' - 0.16 • 4' - 0.65 • 6' • 1.47 • 8' - 2.61 • 10' - 4.08 • 12' - 5.57

Pumpoff 1449

| | amec $^{\odot}$ | LOW-FLOW GROUNDWATER SAMPLING LOG |
|-------|---|---|
| | WELL NO: MW-2 DATE: 3-14-08 OVAPID READING WHEN WELL STATIC WATER LEVEL (TOC): | LOCATION: SPL NOTA TRANSFOR STOTE PROJECT NO: 8915-163410 TIME: 11:36 CLIMATIC CONDITIONS: DVCCCAST LOPENED: NA DEPTH TO PRODUCT (TOC): NA 5.34 TOTAL DEPTH OF WELL (TOC): 4.21 |
| mpon | 1150 METHOD OF I | 1/4 |
| | WELL DTW PURGE DATA | $\begin{array}{c ccccccccccccccccccccccccccccccccccc$ |
| | SAMPLE WITHDRAWAL METHO SAMPLE NUMBER(S) AND TIME | DD: 12.30 7.60 0.658 SAMPLED BY: ME TCH |
| | LAB ANALYSIS PARAMETERS A PESHCI (LO HE) | TENTOMING DODESTY TENDRY MIGH ACTORIGIDA AND PRESERVATIVES: NWITCH-GX, NWITCH-DX, VOC, SVOC, PCB, MICHAELS, LEI DF. HENRY ME FOLD. |
| | NUMBER AND TYPES OF SAMP | The containers used: 5-40 m L 11716, 8-100 m L 91765 |
| | SAMPLES DELIVERED TO: | NSITE TRANSPORTER: |
| HZOCC | DATE: <u>(</u> Diluma = 8.97 1.4 gal Dilu 8.65 | 11ME: 1746 CAPACTY OF CASHIG (CALLOHSUNEAR FOOT) 7.016.11.065.61.11.18.261.10.408.12.557 X3: 4.3 + 6+70 AC PH M3/cm NTU M9/L MU 1223 41/2 7.00 G+6+3 0.1673 ER 141.51 130 12.24 7.00 (COO OPAQUE IT. br) |

| amec $^{\odot}$ | LOW-FI | OW GF | ROUNDV | VATER | SAMPL | ING LO | G | | |
|---|---|---|--|---|---|--|--|--|---------------------|
| WELL NO: MW - 3 DATE 3-14, 08 OVAPID READING WHEN WELL STATIC WATER LEVEL (TOC): | LOCATION: _ TIME: _ OPENED: _ | SPU CA23 13.2 | <u> </u> | CLIMATIC CO DEPTH TO P | STA. ONDITIONS: PRODUCT (TOC TH OF WELL (T | » <u>Iq</u> | 1000 S 1011 104 9,37 | 91516 409 -505 | 03410 Tudlon |
| METHOD OF I | REMOVAL: | 1000- | 90w | Park | HOLDEN | G RATE: | 1/8g/ | min | Γ |
| WELL DTW PURGE 13,83 DATA 14,04 14,09 14,30 14,30 14,48 14,70 14,91 | Time 10 . 14 10 . 16 10 . 18 10 . 20 10 . 24 10 . 26 10 . 28 10 . 30 | Gallons Removed 14 G 10 14 G 14 | 193 3 185 5 181 3 190 5 190 5 190 5 | pH (1) 3.79 (2) 3.71 | 19. Cond. (m8/cm) 2.475 2.489 2.502 2.495 2.506 1.498 2.518 2.514 | Turbidity (NTU) 78.8 36.2 238 14.6 11.2 21.4 9.1 13.9 11.3 | 00 (mg/L) 12.37 11.73 9.88 11.14 11.31 11.64 9.40 11.12 10.5 9.59 | REDOX (mw) 227 218 208 205 201 196 193 199 | did no Calibro |
| SAMPLE WITHDRAWAL METHO SAMPLE NUMBER(S) AND TIME | | 105 | 5 | | | SAMPLED | BY: [ME | - a () 0:57 | <u>'</u> |
| NOTES: O O O | and | ming | prope | slya | on He | SCILLY. | 422 | | |
| LAB ANALYSIS PARAMETERS A NWTPH GIX (H VOOSTU (VON P | | TIVES: | TOX+ | 1+CL | 1225 | 100C 1013 (1 | (mie) |), Hat | E(mc) |
| NUMBER AND TYPES OF SAMP | | S USED: | | 10mL 4, E | 917155 3-1700 | mL C | 3, 2-2 1/2/6 C | Wnles | <u>- glass</u> - |
| DECON. PROCEDURES: SC | ney t | +,0 Cate | ADT C | s l | soles d 10 | level | MHC | 5, 7/ | <u>/</u> |
| SAMPLES DELIVERED TO DATE: | 1514 13.14 | 08 | | TRANSPORT | ER: | 146 | | | |
| HzOcolumn = | HZOCOLUMN = 6.14' CAPACITY OF CASING [GALLONSAINEAR FOOT] 2.018.4.085.6.147.8.261.187.408.12.557 0.988x3 = 2.94 Pump on 10:09 pump off 10:32 | | | | | | | | |
| pumpox | 10:09 | Pu | mp ĉ | P | 10:3 | 2 | | | |

C OnSite Environmental Inc.

Chain of Custody

enutaioM % 03-105 OS.19.08 1746 Metals - Lits 9.452 Chromatograms with final report Requested Analysis 4EM by 1664 ICLP Metals (8) sisteM AROR listo Af6t8 yd sebioidref Pesticides by 8081A CBs by 8082 MIS / GOTS8 vd 2HA9 Laboratory Number: 3/4/1/8 1746 MT2 GOYS8 yd eelitelovimes alogenated Volatiles by 8260B 40928 yd selilalo WTPH-Dx AWTPH-GX/STEX- **UMTPH-HCID** 1 Day 🗌 з Day (TPH analysis 5 working days) 9 2 <u>e_</u> # of Cont. <u>ي</u> \mathcal{U} X Standard (7 working days) Turnaround Request (in working days) 3 Sampled Matrix Reviewed by/Date (Check One) (other) 72 12.F1.1801/C.S.J. Same Day X ☐ 2 Day S.H.C Date Sampled AMÉC EXTRA SAURCOMENTO IMERICASIO CHOCULTURA SPU North Harrie States Phone: (425) 883-3881 • Fax: (425) 885-4603 Sample Identification Project Number:
8-96-165-16541-0 イ・エスとしの YNW) - 2 MW-7 J-90 Reviewed by/Date Relinquished by Relinquished by Retinquished by Received by Received by Received by Lab ID

DISTRIBUTION LEGEND: White - OnSite Copy Yellow - Report Copy Pink - Client Copy



APPENDIX C

Chain-of-Custody Form And Laboratory Analytical Results



14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 20, 2008

Cherilyn Inouye AMEC Earth & Environmental, Inc. 5007 Pacific Highway East, Suite 5 Tacoma, WA 98424

Re: Analytical Data for Project 8915 162890

Laboratory Reference No. 0803-062

Dear Cherilyn:

Enclosed are the analytical results and associated quality control data for samples submitted on March 10, 2008.

Please note that the data for the added Hexavalent Chromium will follow in the final report.

The standard policy of OnSite Environmental Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Project: 8915 162890

Case Narrative

Samples were collected on March 10, 2008 and received by the laboratory on March 10, 2008. They were maintained at the laboratory at a temperature of 2°C to 6°C except as noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

NWTPH Gx Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Volatiles EPA 8260B Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

The result for Toluene in samples B-1_5 and B-3_3 may be the result of contaminated sample vials.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Organochlorine Pesticides by EPA 8081A Analysis

Due to negative effects of the matrix on the instrument, values for the analytes 4,4'-DDT and Methoxychlor in the closing continuing calibration verification standard (CCV) were low. Therefore, values for these compounds can be greater than reported. Since the degradation of the CCV standards was reproducible after re-injecting the sample extracts, the CCV degradation problem was attributed to the matrix of these samples.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Project: 8915 162890

NWTPH-Gx

Date Extracted: 3-12-08
Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Client ID: **B-1_5 B-3_3** Lab ID: 03-062-01 03-062-04

| | Result | Flags | PQL | Result | Flags | PQL |
|--------------------------------------|--------|-------|-----|--------|-------|-----|
| TPH-Gas | ND | | 5.6 | ND | | 6.5 |
| Surrogate Recovery: Fluorobenzene | 94% | | | 94% | | |

Project: 8915 162890

NWTPH-Gx METHOD BLANK QUALITY CONTROL

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0312S1

Result Flags PQL

TPH-Gas ND 5.0

Surrogate Recovery:

Fluorobenzene 99%

Project: 8915 162890

NWTPH-Gx DUPLICATE QUALITY CONTROL

Date Extracted: 3-12-08
Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-01 03-062-01

Original Duplicate RPD Flags

TPH-Gas ND ND NA

Surrogate Recovery:

Fluorobenzene 94% 94%

Project: 8915 162890

NWTPH-Dx

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

 Client ID:
 B-1_5
 B-3_3

 Lab ID:
 03-062-01
 03-062-04

 Diesel Range:
 ND
 150

 PQL:
 29
 31

Identification: --- Diesel Fuel#2

 Lube Oil Range:
 ND
 300

 PQL:
 57
 62

Identification: --- Lube Oil

Surrogate Recovery

o-Terphenyl: 84% 72%

Flags: Y Y

Project: 8915 162890

NWTPH-Dx METHOD BLANK QUALITY CONTROL

| Date Extracted: | 3-12-08 |
|-----------------|---------|
| Date Analyzed: | 3-12-08 |
| | |

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0312S1

Diesel Range: ND PQL: 25

Identification: ---

Lube Oil Range: ND PQL: 50

Identification: ---

Surrogate Recovery

o-Terphenyl: 88%

Flags: Y

Project: 8915 162890

NWTPH-Dx DUPLICATE QUALITY CONTROL

03-062-01 DUP

| Date Extracted: | 3-12-08 |
|-----------------|-------------|
| Date Analyzed: | 3-12-08 |
| Matrix: | Soil |
| Units: | mg/kg (ppm) |

03-062-01

| Diesel Range: | ND | ND |
|---------------|----|----|
| PQL: | 25 | 25 |

RPD: N/A

Surrogate Recovery

Lab ID:

o-Terphenyl: 84% 75%

Flags: Y Y

Project: 8915 162890

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-01 **Client ID: B-1_5**

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|---------|
| Dichlorodifluoromethane | ND | | 0.00090 |
| Chloromethane | ND | | 0.0045 |
| Vinyl Chloride | ND | | 0.00090 |
| Bromomethane | ND | | 0.00090 |
| Chloroethane | ND | | 0.0045 |
| Trichlorofluoromethane | ND | | 0.00090 |
| 1,1-Dichloroethene | ND | | 0.00090 |
| Acetone | 0.014 | | 0.0045 |
| lodomethane | ND | | 0.0045 |
| Carbon Disulfide | ND | | 0.00090 |
| Methylene Chloride | ND | | 0.0045 |
| (trans) 1,2-Dichloroethene | ND | | 0.00090 |
| Methyl t-Butyl Ether | ND | | 0.00090 |
| 1,1-Dichloroethane | ND | | 0.00090 |
| Vinyl Acetate | ND | | 0.0045 |
| 2,2-Dichloropropane | ND | | 0.00090 |
| (cis) 1,2-Dichloroethene | ND | | 0.00090 |
| 2-Butanone | ND | | 0.0045 |
| Bromochloromethane | ND | | 0.00090 |
| Chloroform | ND | | 0.00090 |
| 1,1,1-Trichloroethane | ND | | 0.00090 |
| Carbon Tetrachloride | ND | | 0.00090 |
| 1,1-Dichloropropene | ND | | 0.00090 |
| Benzene | ND | | 0.00090 |
| 1,2-Dichloroethane | ND | | 0.00090 |
| Trichloroethene | ND | | 0.00090 |
| 1,2-Dichloropropane | ND | | 0.00090 |
| Dibromomethane | ND | | 0.00090 |
| Bromodichloromethane | ND | | 0.00090 |
| 2-Chloroethyl Vinyl Ether | ND | | 0.0045 |
| (cis) 1,3-Dichloropropene | ND | | 0.00090 |
| Methyl Isobutyl Ketone | ND | | 0.0045 |
| Toluene | 0.0022 | Z | 0.00090 |
| (trans) 1,3-Dichloropropene | ND | | 0.00090 |

Project: 8915 162890

VOLATILES by EPA 8260B Page 2 of 2

Lab ID: 03-062-01 **Client ID:** B-1_5

| Compound | Results | s Flags | PQL |
|-----------------------------|---------|---------|---------|
| 1,1,2-Trichloroethane | ND | | 0.00090 |
| Tetrachloroethene | ND | | 0.00090 |
| 1,3-Dichloropropane | ND | | 0.00090 |
| 2-Hexanone | ND | | 0.0045 |
| Dibromochloromethane | ND | | 0.00090 |
| 1,2-Dibromoethane | ND | | 0.00090 |
| Chlorobenzene | ND | | 0.00090 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.00090 |
| Ethylbenzene | ND | | 0.00090 |
| m,p-Xylene | ND | | 0.0018 |
| o-Xylene | ND | | 0.00090 |
| Styrene | ND | | 0.00090 |
| Bromoform | ND | | 0.00090 |
| Isopropylbenzene | ND | | 0.00090 |
| Bromobenzene | ND | | 0.00090 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.00090 |
| 1,2,3-Trichloropropane | ND | | 0.00090 |
| n-Propylbenzene | ND | | 0.00090 |
| 2-Chlorotoluene | ND | | 0.00090 |
| 4-Chlorotoluene | ND | | 0.00090 |
| 1,3,5-Trimethylbenzene | ND | | 0.00090 |
| tert-Butylbenzene | ND | | 0.00090 |
| 1,2,4-Trimethylbenzene | ND | | 0.00090 |
| sec-Butylbenzene | ND | | 0.00090 |
| 1,3-Dichlorobenzene | ND | | 0.00090 |
| p-Isopropyltoluene | ND | | 0.00090 |
| 1,4-Dichlorobenzene | ND | | 0.00090 |
| 1,2-Dichlorobenzene | ND | | 0.00090 |
| n-Butylbenzene | ND | | 0.00090 |
| 1,2-Dibromo-3-chloropropane | ND | | 0.0045 |
| 1,2,4-Trichlorobenzene | ND | | 0.00090 |
| Hexachlorobutadiene | ND | | 0.0045 |
| Naphthalene | ND | | 0.00090 |
| 1,2,3-Trichlorobenzene | ND | | 0.00090 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 77 | 70-118 |
| Toluene-d8 | 95 | 70-121 |
| 4-Bromofluorobenzene | 82 | 70-130 |

Project: 8915 162890

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-04 **Client ID: B-3_3**

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|--------|
| Dichlorodifluoromethane | ND | | 0.0011 |
| Chloromethane | ND | | 0.0053 |
| Vinyl Chloride | ND | | 0.0011 |
| Bromomethane | ND | | 0.0011 |
| Chloroethane | ND | | 0.0053 |
| Trichlorofluoromethane | ND | | 0.0011 |
| 1,1-Dichloroethene | ND | | 0.0011 |
| Acetone | ND | | 0.0053 |
| lodomethane | ND | | 0.0053 |
| Carbon Disulfide | ND | | 0.0011 |
| Methylene Chloride | ND | | 0.0053 |
| (trans) 1,2-Dichloroethene | ND | | 0.0011 |
| Methyl t-Butyl Ether | ND | | 0.0011 |
| 1,1-Dichloroethane | ND | | 0.0011 |
| Vinyl Acetate | ND | | 0.0053 |
| 2,2-Dichloropropane | ND | | 0.0011 |
| (cis) 1,2-Dichloroethene | ND | | 0.0011 |
| 2-Butanone | ND | | 0.0053 |
| Bromochloromethane | ND | | 0.0011 |
| Chloroform | ND | | 0.0011 |
| 1,1,1-Trichloroethane | ND | | 0.0011 |
| Carbon Tetrachloride | ND | | 0.0011 |
| 1,1-Dichloropropene | ND | | 0.0011 |
| Benzene | ND | | 0.0011 |
| 1,2-Dichloroethane | ND | | 0.0011 |
| Trichloroethene | ND | | 0.0011 |
| 1,2-Dichloropropane | ND | | 0.0011 |
| Dibromomethane | ND | | 0.0011 |
| Bromodichloromethane | ND | | 0.0011 |
| 2-Chloroethyl Vinyl Ether | ND | | 0.0053 |
| (cis) 1,3-Dichloropropene | ND | | 0.0011 |
| Methyl Isobutyl Ketone | ND | | 0.0053 |
| Toluene | 0.0030 | Z | 0.0011 |
| (trans) 1,3-Dichloropropene | ND | | 0.0011 |

Project: 8915 162890

VOLATILES by EPA 8260B Page 2 of 2

Lab ID: 03-062-04 **Client ID:** B-3_3

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|--------|
| 1,1,2-Trichloroethane | ND | _ | 0.0011 |
| Tetrachloroethene | ND | | 0.0011 |
| 1,3-Dichloropropane | ND | | 0.0011 |
| 2-Hexanone | ND | | 0.0053 |
| Dibromochloromethane | ND | | 0.0011 |
| 1,2-Dibromoethane | ND | | 0.0011 |
| Chlorobenzene | ND | | 0.0011 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.0011 |
| Ethylbenzene | ND | | 0.0011 |
| m,p-Xylene | ND | | 0.0021 |
| o-Xylene | ND | | 0.0011 |
| Styrene | ND | | 0.0011 |
| Bromoform | ND | | 0.0011 |
| Isopropylbenzene | ND | | 0.0011 |
| Bromobenzene | ND | | 0.0011 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.0011 |
| 1,2,3-Trichloropropane | ND | | 0.0011 |
| n-Propylbenzene | ND | | 0.0011 |
| 2-Chlorotoluene | ND | | 0.0011 |
| 4-Chlorotoluene | ND | | 0.0011 |
| 1,3,5-Trimethylbenzene | ND | | 0.0011 |
| tert-Butylbenzene | ND | | 0.0011 |
| 1,2,4-Trimethylbenzene | ND | | 0.0011 |
| sec-Butylbenzene | ND | | 0.0011 |
| 1,3-Dichlorobenzene | ND | | 0.0011 |
| p-Isopropyltoluene | ND | | 0.0011 |
| 1,4-Dichlorobenzene | ND | | 0.0011 |
| 1,2-Dichlorobenzene | ND | | 0.0011 |
| n-Butylbenzene | ND | | 0.0011 |
| 1,2-Dibromo-3-chloropropane | ND | | 0.0053 |
| 1,2,4-Trichlorobenzene | ND | | 0.0011 |
| Hexachlorobutadiene | ND | | 0.0053 |
| Naphthalene | ND | | 0.0011 |
| 1,2,3-Trichlorobenzene | ND | | 0.0011 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 82 | 70-118 |
| Toluene-d8 | 90 | 70-121 |
| 4-Bromofluorobenzene | 95 | 70-130 |

Project: 8915 162890

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 1 of 2

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0314S1

| Paculte | Flage | PQL |
|---------|--|--|
| | i lags | 0.0010 |
| | | 0.0050 |
| | | 0.0010 |
| | | 0.0010 |
| | | 0.0050 |
| | | 0.0010 |
| | | 0.0010 |
| ND | | 0.0050 |
| ND | | 0.0050 |
| ND | | 0.0010 |
| ND | | 0.0050 |
| ND | | 0.0010 |
| ND | | 0.0010 |
| ND | | 0.0010 |
| ND | | 0.0050 |
| ND | | 0.0010 |
| ND | | 0.0010 |
| ND | | 0.0050 |
| ND | | 0.0010 |
| ND | | 0.0050 |
| | | 0.0010 |
| ND | | 0.0050 |
| ND | | 0.0010 |
| ND | | 0.0010 |
| | ND N | ND N |

Project: 8915 162890

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 2 of 2

Lab ID: MB0314S1

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|---------|
| 1,1,2-Trichloroethane | ND | | 0.0010 |
| Tetrachloroethene | ND | | 0.0010 |
| 1,3-Dichloropropane | ND | | 0.0010 |
| 2-Hexanone | ND | | 0.0050 |
| Dibromochloromethane | ND | | 0.0010 |
| 1,2-Dibromoethane | ND | | 0.0010 |
| Chlorobenzene | ND | | 0.0010 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.0010 |
| Ethylbenzene | ND | | 0.0010 |
| m,p-Xylene | ND | | 0.0020 |
| o-Xylene | ND | | 0.0010 |
| Styrene | ND | | 0.0010 |
| Bromoform | ND | | 0.0010 |
| Isopropylbenzene | ND | | 0.0010 |
| Bromobenzene | ND | | 0.0010 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.0010 |
| 1,2,3-Trichloropropane | ND | | 0.0010 |
| n-Propylbenzene | ND | | 0.0010 |
| 2-Chlorotoluene | ND | | 0.0010 |
| 4-Chlorotoluene | ND | | 0.0010 |
| 1,3,5-Trimethylbenzene | ND | | 0.0010 |
| tert-Butylbenzene | ND | | 0.0010 |
| 1,2,4-Trimethylbenzene | ND | | 0.0010 |
| sec-Butylbenzene | ND | | 0.0010 |
| 1,3-Dichlorobenzene | ND | | 0.0010 |
| p-Isopropyltoluene | ND | | 0.0010 |
| 1,4-Dichlorobenzene | ND | | 0.0010 |
| 1,2-Dichlorobenzene | ND | | 0.0010 |
| n-Butylbenzene | ND | | 0.0010 |
| 1,2-Dibromo-3-chloropropane | ND | | 0.0050 |
| 1,2,4-Trichlorobenzene | ND | | 0.0010 |
| Hexachlorobutadiene | ND | | 0.0050 |
| Naphthalene | ND | | 0.0010 |
| 1,2,3-Trichlorobenzene | ND | | 0.0010 |
| Surregate | Percent | | Control |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 70 | 70-118 |
| Toluene-d8 | 91 | 70-121 |
| 4-Bromofluorobenzene | 79 | 70-130 |

Project: 8915 162890

VOLATILES by EPA 8260B SB/SBD QUALITY CONTROL

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: SB0314S1

| | Spike | | Percent | | Percent | Recovery | |
|--------------------|--------|--------|----------|--------|----------|----------|-------|
| Compound | Amount | SB | Recovery | SBD | Recovery | Limits | Flags |
| | | | | | | | |
| 1,1-Dichloroethene | 0.0500 | 0.0373 | 75 | 0.0412 | 82 | 70-130 | |
| Benzene | 0.0500 | 0.0419 | 84 | 0.0422 | 84 | 70-127 | |
| Trichloroethene | 0.0500 | 0.0404 | 81 | 0.0449 | 90 | 73-117 | |
| Toluene | 0.0500 | 0.0410 | 82 | 0.0436 | 87 | 78-115 | |
| Chlorobenzene | 0.0500 | 0.0410 | 82 | 0.0418 | 84 | 80-117 | |

| | RPD | | |
|--------------------|-----|-------|-------|
| | RPD | Limit | Flags |
| 1,1-Dichloroethene | 10 | 10 | |
| Benzene | 1 | 11 | |
| Trichloroethene | 11 | 13 | |
| Toluene | 6 | 12 | |
| Chlorobenzene | 2 | 10 | |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 1 of 3

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-03 **Client ID: B-1_8**

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|--------|
| N-Nitrosodimethylamine | ND | _ | 0.037 |
| Pyridine | ND | | 0.037 |
| Phenol | ND | | 0.037 |
| Aniline | ND | | 0.037 |
| bis(2-Chloroethyl)ether | ND | | 0.037 |
| 2-Chlorophenol | ND | | 0.037 |
| 1,3-Dichlorobenzene | ND | | 0.037 |
| 1,4-Dichlorobenzene | ND | | 0.037 |
| Benzyl alcohol | ND | | 0.037 |
| 1,2-Dichlorobenzene | ND | | 0.037 |
| 2-Methylphenol (o-Cresol) | ND | | 0.037 |
| bis(2-Chloroisopropyl)ether | ND | | 0.037 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 0.037 |
| N-Nitroso-di-n-propylamine | ND | | 0.037 |
| Hexachloroethane | ND | | 0.037 |
| Nitrobenzene | ND | | 0.037 |
| Isophorone | ND | | 0.037 |
| 2-Nitrophenol | ND | | 0.037 |
| 2,4-Dimethylphenol | ND | | 0.037 |
| bis(2-Chloroethoxy)methane | ND | | 0.037 |
| 2,4-Dichlorophenol | ND | | 0.037 |
| 1,2,4-Trichlorobenzene | ND | | 0.037 |
| Naphthalene | ND | | 0.0074 |
| 4-Chloroaniline | ND | | 0.037 |
| Hexachlorobutadiene | ND | | 0.037 |
| 4-Chloro-3-methylphenol | ND | | 0.037 |
| 2-Methylnaphthalene | ND | | 0.0074 |
| 1-Methylnaphthalene | ND | | 0.0074 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

Lab ID: 03-062-03 **Client ID: B-1_8**

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Hexachlorocyclopentadiene | ND | | 0.037 |
| 2,4,6-Trichlorophenol | ND | | 0.037 |
| 2,3-Dichloroaniline | ND | | 0.037 |
| 2,4,5-Trichlorophenol | ND | | 0.037 |
| 2-Chloronaphthalene | ND | | 0.037 |
| 2-Nitroaniline | ND | | 0.037 |
| 1,4-Dinitrobenzene | ND | | 0.037 |
| Dimethylphthalate | ND | | 0.037 |
| 1,3-Dinitrobenzene | ND | | 0.037 |
| 2,6-Dinitrotoluene | ND | | 0.037 |
| 1,2-Dinitrobenzene | ND | | 0.037 |
| Acenaphthylene | ND | | 0.0074 |
| 3-Nitroaniline | ND | | 0.037 |
| 2,4-Dinitrophenol | ND | | 0.19 |
| Acenaphthene | ND | | 0.0074 |
| 4-Nitrophenol | ND | | 0.037 |
| 2,4-Dinitrotoluene | ND | | 0.037 |
| Dibenzofuran | ND | | 0.037 |
| 2,3,4,6-Tetrachlorophenol | ND | | 0.037 |
| 2,3,5,6-Tetrachlorophenol | ND | | 0.037 |
| Diethylphthalate | ND | | 0.037 |
| 4-Chlorophenyl-phenylether | ND | | 0.037 |
| 4-Nitroaniline | ND | | 0.037 |
| Fluorene | ND | | 0.0074 |
| 4,6-Dinitro-2-methylphenol | ND | | 0.19 |
| N-Nitrosodiphenylamine | ND | | 0.037 |
| 1,2-Diphenylhydrazine | ND | | 0.037 |
| 4-Bromophenyl-phenylether | ND | | 0.037 |
| Hexachlorobenzene | ND | | 0.037 |
| Pentachlorophenol | ND | | 0.19 |
| Phenanthrene | ND | | 0.0074 |
| Anthracene | ND | | 0.0074 |
| Carbazole | ND | | 0.037 |
| Di-n-butylphthalate | 0.041 | | 0.037 |
| Fluoranthene | ND | | 0.0074 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 3 of 3

Lab ID: 03-062-03 **Client ID: B-1_8**

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Benzidine | ND | _ | 0.37 |
| Pyrene | ND | | 0.0074 |
| Butylbenzylphthalate | ND | | 0.037 |
| bis-2-Ethylhexyladipate | ND | | 0.037 |
| 3,3'-Dichlorobenzidine | ND | | 0.037 |
| Benzo[a]anthracene | ND | | 0.0074 |
| Chrysene | ND | | 0.0074 |
| bis(2-Ethylhexyl)phthalate | ND | | 0.037 |
| Di-n-octylphthalate | ND | | 0.037 |
| Benzo[b]fluoranthene | ND | | 0.0074 |
| Benzo[k]fluoranthene | ND | | 0.0074 |
| Benzo[a]pyrene | ND | | 0.0074 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.0074 |
| Dibenz[a,h]anthracene | ND | | 0.0074 |
| Benzo[g,h,i]perylene | ND | | 0.0074 |

| Surrogate : | ogate : Percent Recovery | | | |
|----------------------|--------------------------|----------|--|--|
| | | | | |
| 2-Fluorophenol | 73 | 39 - 90 | | |
| Phenol-d6 | 88 | 40 - 100 | | |
| Nitrobenzene-d5 | 81 | 30 - 100 | | |
| 2-Fluorobiphenyl | 84 | 41 - 100 | | |
| 2,4,6-Tribromophenol | 88 | 53 - 105 | | |
| Terphenyl-d14 | 100 | 49 - 115 | | |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 1 of 3

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-04 **Client ID: B-3_3**

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|--------|
| N-Nitrosodimethylamine | ND | • | 0.041 |
| Pyridine | ND | | 0.041 |
| Phenol | ND | | 0.041 |
| Aniline | ND | | 0.041 |
| bis(2-Chloroethyl)ether | ND | | 0.041 |
| 2-Chlorophenol | ND | | 0.041 |
| 1,3-Dichlorobenzene | ND | | 0.041 |
| 1,4-Dichlorobenzene | ND | | 0.041 |
| Benzyl alcohol | ND | | 0.041 |
| 1,2-Dichlorobenzene | ND | | 0.041 |
| 2-Methylphenol (o-Cresol) | ND | | 0.041 |
| bis(2-Chloroisopropyl)ether | ND | | 0.041 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 0.041 |
| N-Nitroso-di-n-propylamine | ND | | 0.041 |
| Hexachloroethane | ND | | 0.041 |
| Nitrobenzene | ND | | 0.041 |
| Isophorone | ND | | 0.041 |
| 2-Nitrophenol | ND | | 0.041 |
| 2,4-Dimethylphenol | ND | | 0.041 |
| bis(2-Chloroethoxy)methane | ND | | 0.041 |
| 2,4-Dichlorophenol | ND | | 0.041 |
| 1,2,4-Trichlorobenzene | ND | | 0.041 |
| Naphthalene | ND | | 0.0082 |
| 4-Chloroaniline | ND | | 0.041 |
| Hexachlorobutadiene | ND | | 0.041 |
| 4-Chloro-3-methylphenol | ND | | 0.041 |
| 2-Methylnaphthalene | ND | | 0.0082 |
| 1-Methylnaphthalene | ND | | 0.0082 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

Lab ID: 03-062-04 **Client ID: B-3_3**

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Hexachlorocyclopentadiene | ND | • | 0.041 |
| 2,4,6-Trichlorophenol | ND | | 0.041 |
| 2,3-Dichloroaniline | ND | | 0.041 |
| 2,4,5-Trichlorophenol | ND | | 0.041 |
| 2-Chloronaphthalene | ND | | 0.041 |
| 2-Nitroaniline | ND | | 0.041 |
| 1,4-Dinitrobenzene | ND | | 0.041 |
| Dimethylphthalate | ND | | 0.041 |
| 1,3-Dinitrobenzene | ND | | 0.041 |
| 2,6-Dinitrotoluene | ND | | 0.041 |
| 1,2-Dinitrobenzene | ND | | 0.041 |
| Acenaphthylene | ND | | 0.0082 |
| 3-Nitroaniline | ND | | 0.041 |
| 2,4-Dinitrophenol | ND | | 0.21 |
| Acenaphthene | ND | | 0.0082 |
| 4-Nitrophenol | ND | | 0.041 |
| 2,4-Dinitrotoluene | ND | | 0.041 |
| Dibenzofuran | ND | | 0.041 |
| 2,3,4,6-Tetrachlorophenol | ND | | 0.041 |
| 2,3,5,6-Tetrachlorophenol | ND | | 0.041 |
| Diethylphthalate | ND | | 0.041 |
| 4-Chlorophenyl-phenylether | ND | | 0.041 |
| 4-Nitroaniline | ND | | 0.041 |
| Fluorene | ND | | 0.0082 |
| 4,6-Dinitro-2-methylphenol | ND | | 0.21 |
| N-Nitrosodiphenylamine | ND | | 0.041 |
| 1,2-Diphenylhydrazine | ND | | 0.041 |
| 4-Bromophenyl-phenylether | ND | | 0.041 |
| Hexachlorobenzene | ND | | 0.041 |
| Pentachlorophenol | ND | | 0.21 |
| Phenanthrene | 0.013 | | 0.0082 |
| Anthracene | ND | | 0.0082 |
| Carbazole | ND | | 0.041 |
| Di-n-butylphthalate | 0.071 | | 0.041 |
| Fluoranthene | ND | | 0.0082 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 3 of 3

Lab ID: 03-062-04 **Client ID: B-3_3**

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Benzidine | ND | | 0.41 |
| Pyrene | 0.011 | | 0.0082 |
| Butylbenzylphthalate | ND | | 0.041 |
| bis-2-Ethylhexyladipate | ND | | 0.041 |
| 3,3'-Dichlorobenzidine | ND | | 0.041 |
| Benzo[a]anthracene | ND | | 0.0082 |
| Chrysene | ND | | 0.0082 |
| bis(2-Ethylhexyl)phthalate | 0.052 | | 0.041 |
| Di-n-octylphthalate | ND | | 0.041 |
| Benzo[b]fluoranthene | ND | | 0.0082 |
| Benzo[k]fluoranthene | ND | | 0.0082 |
| Benzo[a]pyrene | ND | | 0.0082 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.0082 |
| Dibenz[a,h]anthracene | ND | | 0.0082 |
| Benzo[g,h,i]perylene | ND | | 0.0082 |

| Surrogate : | Percent Recovery | |
|----------------------|---------------------|----------|
| 2-Fluorophenol | 77 | 39 - 90 |
| Phenol-d6 | 94 | 40 - 100 |
| Nitrobenzene-d5 | 85 | 30 - 100 |
| 2-Fluorobiphenyl | 88 | 41 - 100 |
| 2,4,6-Tribromophenol | 91 | 53 - 105 |
| Terphenyl-d14 | 98 | 49 - 115 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 1 of 3

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0312S1

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|--------|
| N-Nitrosodimethylamine | ND | | 0.033 |
| Pyridine | ND | | 0.033 |
| Phenol | ND | | 0.033 |
| Aniline | ND | | 0.033 |
| bis(2-Chloroethyl)ether | ND | | 0.033 |
| 2-Chlorophenol | ND | | 0.033 |
| 1,3-Dichlorobenzene | ND | | 0.033 |
| 1,4-Dichlorobenzene | ND | | 0.033 |
| Benzyl alcohol | ND | | 0.033 |
| 1,2-Dichlorobenzene | ND | | 0.033 |
| 2-Methylphenol (o-Cresol) | ND | | 0.033 |
| bis(2-Chloroisopropyl)ether | ND | | 0.033 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 0.033 |
| N-Nitroso-di-n-propylamine | ND | | 0.033 |
| Hexachloroethane | ND | | 0.033 |
| Nitrobenzene | ND | | 0.033 |
| Isophorone | ND | | 0.033 |
| 2-Nitrophenol | ND | | 0.033 |
| 2,4-Dimethylphenol | ND | | 0.033 |
| bis(2-Chloroethoxy)methane | ND | | 0.033 |
| 2,4-Dichlorophenol | ND | | 0.033 |
| 1,2,4-Trichlorobenzene | ND | | 0.033 |
| Naphthalene | ND | | 0.0067 |
| 4-Chloroaniline | ND | | 0.033 |
| Hexachlorobutadiene | ND | | 0.033 |
| 4-Chloro-3-methylphenol | ND | | 0.033 |
| 2-Methylnaphthalene | ND | | 0.0067 |
| 1-Methylnaphthalene | ND | | 0.0067 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 2 of 3

Lab ID: MB0312S1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Hexachlorocyclopentadiene | ND | · · | 0.033 |
| 2,4,6-Trichlorophenol | ND | | 0.033 |
| 2,3-Dichloroaniline | ND | | 0.033 |
| 2,4,5-Trichlorophenol | ND | | 0.033 |
| 2-Chloronaphthalene | ND | | 0.033 |
| 2-Nitroaniline | ND | | 0.033 |
| 1,4-Dinitrobenzene | ND | | 0.033 |
| Dimethylphthalate | ND | | 0.033 |
| 1,3-Dinitrobenzene | ND | | 0.033 |
| 2,6-Dinitrotoluene | ND | | 0.033 |
| 1,2-Dinitrobenzene | ND | | 0.033 |
| Acenaphthylene | ND | | 0.0067 |
| 3-Nitroaniline | ND | | 0.033 |
| 2,4-Dinitrophenol | ND | | 0.17 |
| Acenaphthene | ND | | 0.0067 |
| 4-Nitrophenol | ND | | 0.033 |
| 2,4-Dinitrotoluene | ND | | 0.033 |
| Dibenzofuran | ND | | 0.033 |
| 2,3,4,6-Tetrachlorophenol | ND | | 0.033 |
| 2,3,5,6-Tetrachlorophenol | ND | | 0.033 |
| Diethylphthalate | ND | | 0.033 |
| 4-Chlorophenyl-phenylether | ND | | 0.033 |
| 4-Nitroaniline | ND | | 0.033 |
| Fluorene | ND | | 0.0067 |
| 4,6-Dinitro-2-methylphenol | ND | | 0.17 |
| N-Nitrosodiphenylamine | ND | | 0.033 |
| 1,2-Diphenylhydrazine | ND | | 0.033 |
| 4-Bromophenyl-phenylether | ND | | 0.033 |
| Hexachlorobenzene | ND | | 0.033 |
| Pentachlorophenol | ND | | 0.17 |
| Phenanthrene | ND | | 0.0067 |
| Anthracene | ND | | 0.0067 |
| Carbazole | ND | | 0.033 |
| Di-n-butylphthalate | ND | | 0.033 |
| Fluoranthene | ND | | 0.0067 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 3 of 3

Lab ID: MB0312S1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Benzidine | ND | | 0.33 |
| Pyrene | ND | | 0.0067 |
| Butylbenzylphthalate | ND | | 0.033 |
| bis-2-Ethylhexyladipate | ND | | 0.033 |
| 3,3'-Dichlorobenzidine | ND | | 0.033 |
| Benzo[a]anthracene | ND | | 0.0067 |
| Chrysene | ND | | 0.0067 |
| bis(2-Ethylhexyl)phthalate | ND | | 0.033 |
| Di-n-octylphthalate | ND | | 0.033 |
| Benzo[b]fluoranthene | ND | | 0.0067 |
| Benzo[k]fluoranthene | ND | | 0.0067 |
| Benzo[a]pyrene | ND | | 0.0067 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.0067 |
| Dibenz[a,h]anthracene | ND | | 0.0067 |
| Benzo[g,h,i]perylene | ND | | 0.0067 |

| Surrogate : | Percent Recovery | Control Limits |
|----------------------|---------------------|-------------------|
| 2-Fluorophenol | 76 | 39 - 90 |
| Phenol-d6 | 90 | 40 - 100 |
| Nitrobenzene-d5 | 84 | 30 - 100 |
| 2-Fluorobiphenyl | 87 | 41 - 100 |
| 2,4,6-Tribromophenol | 90 | 53 - 105 |
| Terphenyl-d14 | 103 | 49 - 115 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM SB/SBD QUALITY CONTROL

Date Extracted: 3-12-08
Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: SB0312S1

| | Spike | | Percent | | Percent | Recovery | , |
|----------------------------|--------|-------|----------|-------|----------|----------|-------|
| Compound: | Amount | SB | Recovery | SBD | Recovery | Limits | Flags |
| Phenol | 1.33 | 1.12 | 84 | 0.993 | 75 | 36-97 | |
| 2-Chlorophenol | 1.33 | 1.10 | 83 | 0.950 | 71 | 32-100 | |
| 1,4-Dichlorobenzene | 0.667 | 0.450 | 67 | 0.382 | 57 | 24-94 | |
| N-Nitroso-di-n-propylamine | 0.667 | 0.567 | 85 | 0.483 | 72 | 34-99 | |
| 1,2,4-Trichlorobenzene | 0.667 | 0.429 | 64 | 0.368 | 55 | 23-85 | |
| 4-Chloro-3-methylphenol | 1.33 | 1.18 | 89 | 1.18 | 88 | 46-108 | |
| Acenaphthene | 0.667 | 0.527 | 79 | 0.499 | 75 | 37-101 | |
| 2,4-Dinitrotoluene | 0.667 | 0.605 | 91 | 0.606 | 91 | 41-116 | |
| 4-Nitrophenol | 1.33 | 1.44 | 108 | 1.50 | 113 | 48-116 | |
| Pentachlorophenol | 1.33 | 1.21 | 91 | 1.25 | 94 | 28-130 | |
| Pyrene | 0.667 | 0.631 | 95 | 0.649 | 97 | 46-117 | |

| | | RPD | |
|----------------------------|-----|--------|-------|
| | RPD | Limits | Flags |
| Phenol | 12 | 29 | |
| 2-Chlorophenol | 14 | 28 | |
| 1,4-Dichlorobenzene | 16 | 27 | |
| N-Nitroso-di-n-propylamine | 16 | 30 | |
| 1,2,4-Trichlorobenzene | 15 | 28 | |
| 4-Chloro-3-methylphenol | 0 | 22 | |
| Acenaphthene | 6 | 25 | |
| 2,4-Dinitrotoluene | 0 | 30 | |
| 4-Nitrophenol | 4 | 30 | |
| Pentachlorophenol | 3 | 30 | |
| Pyrene | 3 | 20 | |
| | | | |

Project: 8915 162890

PCBs by EPA 8082

Matrix: Soil

Units: mg/Kg (ppm)

| 5 5 41 7 | | | | Date | Date | |
|----------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | B-1_3 | | | | | |
| Laboratory ID: | 03-062-02 | | | | | |
| Aroclor 1016 | ND | 0.057 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1221 | ND | 0.057 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1232 | ND | 0.057 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1242 | ND | 0.057 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1248 | ND | 0.057 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1254 | ND | 0.057 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1260 | ND | 0.057 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1262 | ND | 0.057 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1268 | ND | 0.057 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| DCB | 85 | 39-118 | | | | |
| Client ID: | B-3_6 | | | | | |
| Laboratory ID: | 03-062-05 | | | | | |
| Aroclor 1016 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1221 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1232 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1242 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1248 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1254 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1260 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1262 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1268 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |

Surrogate: Percent Recovery Control Limit
DCB 90 39-118

Project: 8915 162890

PCBs by EPA 8082 QUALITY CONTROL

Matrix: Soil

Units: mg/Kg (ppm)

| | | | | Date | Date | |
|----------------|----------|-------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| METHOD BLANK | | | | | | |
| Laboratory ID: | MB0312S1 | | | | | |
| Aroclor 1016 | ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1221 | ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1232 | ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1242 | ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1248 | ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1254 | ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1260 | ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1262 | ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1268 | ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| • | | | | | • | • |

Surrogate: Percent Recovery Control Limits DCB 101 39-118

| Analyte | Re | sult | Spike | Level | Source Result | | rcent covery | Recovery Limits | RPD | RPD Limit | Flags |
|----------------|-------|-------|-------|-------|------------------|----|-----------------|--------------------|-----|--------------|-------|
| MATRIX SPIKES | | | | | | | | | | | |
| Laboratory ID: | 03-0 | 62-05 | | | | | | | | | |
| | MS | MSD | MS | MSD | | MS | MSD | | | | |
| Aroclor 1260 | 0.444 | 0.456 | 0.500 | 0.500 | ND | 89 | 91 | 35-120 | 3 | 17 | |
| Surrogate: | | | | | | | | | | | |
| DCB | | | | | | 93 | 92 | 39-118 | | | |

Project: 8915 162890

ORGANOCHLORINE PESTICIDES by EPA 8081A

Matrix: Soil

Units: ug/Kg (ppb)

| | | | | Date | Date | |
|--------------------|-----------|-----|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | B-1_3 | | | | | |
| Laboratory ID: | 03-062-02 | | | | | |
| alpha-BHC | ND | 5.7 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-BHC | ND | 5.7 | EPA 8081 | 3-12-08 | 3-12-08 | |
| beta-BHC | ND | 5.7 | EPA 8081 | 3-12-08 | 3-12-08 | |
| delta-BHC | ND | 5.7 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor | ND | 5.7 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Aldrin | ND | 5.7 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor Epoxide | ND | 5.7 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-Chlordane | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| alpha-Chlordane | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDE | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan I | ND | 5.7 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Dieldrin | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDD | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan II | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDT | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Aldehyde | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Methoxychlor | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endsulfan Sulfate | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Ketone | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Toxaphene | ND | 57 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 0 | D D | 0 | | · | · | |

Surrogate: Percent Recovery Control Limits TCMX 72 36-108 DCB 63 30-115

Project: 8915 162890

ORGANOCHLORINE PESTICIDES by EPA 8081A

Matrix: Soil

Units: ug/Kg (ppb)

| | | | | Date | Date | |
|--------------------|-----------|-----|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | B-3_6 | | | | | |
| Laboratory ID: | 03-062-05 | | | | | |
| alpha-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| beta-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| delta-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Aldrin | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor Epoxide | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-Chlordane | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| alpha-Chlordane | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDE | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan I | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Dieldrin | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDD | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan II | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDT | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Aldehyde | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Methoxychlor | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endsulfan Sulfate | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Ketone | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Toxaphene | ND | 50 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 0 | D D | 0 | | | | |

Surrogate: Percent Recovery Control Limits TCMX 78 36-108 DCB 75 30-115

Project: 8915 162890

ORGANOCHLORINE PESTICIDES by EPA 8081A QUALITY CONTROL

Matrix: Soil

Units: ug/Kg (ppb)

| Analyte | Result | PQL | Method | Date Propaged | Date Analyzed | Flags |
|--------------------|------------------|----------------|----------|------------------|------------------|-------|
| METHOD BLANK | Result | FQL | Metriou | Prepared | Allalyzeu | riays |
| Laboratory ID: | MB0312S1 | | | | | |
| alpha-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| beta-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| delta-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Aldrin | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor Epoxide | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-Chlordane | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| alpha-Chlordane | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDE | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan I | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Dieldrin | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDD | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan II | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDT | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Aldehyde | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Methoxychlor | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endsulfan Sulfate | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Ketone | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Toxaphene | ND | 50 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| TCMX | 82 | 36-108 | | | | |

Surrogate: Percent Recovery Control Limi
TCMX 82 36-108
DCB 77 30-115

| Analyte | Res | sult | Spike | Level | Source Result | | rcent covery | Recovery Limits | RPD | RPD Limit | Flags |
|----------------|------|-------|-------|-------|------------------|----|-----------------|--------------------|-----|--------------|-------|
| MATRIX SPIKES | | | • | | | | | | | | |
| Laboratory ID: | 03-0 | 62-05 | | | | | | | | | |
| | MS | MSD | MS | MSD | | MS | MSD | | | | |
| gamma-BHC | 41.8 | 39.3 | 50.0 | 50.0 | ND | 84 | 79 | 39-106 | 6 | 12 | |
| Heptachlor | 39.3 | 37.3 | 50.0 | 50.0 | ND | 79 | 75 | 33-107 | 5 | 13 | |
| Aldrin | 39.2 | 37.4 | 50.0 | 50.0 | ND | 78 | 75 | 36-101 | 5 | 12 | |
| Dieldrin | 98.1 | 93.0 | 125 | 125 | ND | 78 | 74 | 33-115 | 5 | 11 | |
| Endrin | 99.4 | 94.8 | 125 | 125 | ND | 80 | 76 | 35-108 | 5 | 11 | |
| 4,4'-DDT | 98.4 | 94.3 | 125 | 125 | ND | 79 | 75 | 24-122 | 4 | 17 | |
| Surrogate: | | | | | | | | | | | |
| TCMX | | | | | | 74 | 71 | 36-108 | | | |
| DCB | | | | | | 72 | 70 | 30-115 | | | |

OnSite Environmental, Inc. 14648 NE 95th Street, Redmond, WA 98052 (425) 883-3881

Project: 8915 162890

CHLORINATED ACID HERBICIDES by EPA 8151A

Matrix: Soil

2,4,5-T

2,4-DB

Dinoseb

Units: ug/Kg (ppb)

| | | | | Date | Date | |
|-------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | B-1_3 | | | | | |
| Laboratory ID: | 03-062-02 | | | | | |
| Dalapon | ND | 260 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 5300 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 5300 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 54 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 1.1 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 54 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-T | ND | 54 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-DB | ND | 54 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dinoseb | ND | 54 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| DCAA | 72 | 37-114 | | | | |
| Client ID: | B-3_6 | | | | | |
| Laboratory ID: | 03-062-05 | | | | | |
| Dalapon | ND | 250 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 52 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 5200 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 5200 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 52 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 52 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 1.1 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| · · · | | | | | | |

Surrogate: Percent Recovery Control Limits DCAA 68 37-114

ND

ND

ND

53

53

53

EPA 8151

EPA 8151

EPA 8151

3-18-08

3-18-08

3-18-08

3-18-08

3-18-08

3-18-08

Project: 8915 162890

CHLORINATED ACID HERBICIDES by EPA 8151A QUALITY CONTROL

Matrix: Soil

Units: ug/Kg (ppb)

| | | | | Date | Date | |
|-------------------|----------|------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| METHOD BLANK | | | | | | |
| Laboratory ID: | MB0318S1 | | | | | |
| Dalapon | ND | 230 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 4700 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 4700 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 0.95 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 48 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-T | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-DB | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dinoseb | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |

Surrogate: Percent Recovery Control Limits DCAA 70 37-114

| | | | | | Source | Pe | rcent | Recovery | | RPD | |
|----------------|------|-------|-------|-------|--------|-----|-------|----------|-----|-------|-------|
| Analyte | Re | sult | Spike | Level | Result | Red | overy | Limits | RPD | Limit | Flags |
| MATRIX SPIKES | | | | | | | | | | | |
| Laboratory ID: | 03-0 | 62-05 | | | | | | | | | |
| | MS | MSD | MS | MSD | | MS | MSD | | | | |
| Dicamba | 77.7 | 76.9 | 100 | 100 | ND | 78 | 77 | 62-99 | 1 | 10 | |
| 2,4-D | 71.5 | 69.1 | 100 | 100 | ND | 72 | 69 | 23-111 | 3 | 24 | |
| 2,4,5-T | 79.5 | 75.5 | 100 | 100 | ND | 79 | 76 | 38-120 | 5 | 12 | |
| 2,4-DB | 90.3 | 83.9 | 100 | 100 | ND | 90 | 84 | 44-135 | 7 | 16 | |
| Surrogate: | | | | | | | | | | | |
| DCAA | | | | | | 80 | 79 | 37-114 | | | |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A

Date Extracted: 3-13&14-08 Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-02 **Client ID: B-1_3**

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 6010B | ND | 11 |
| Barium | 6010B | 46 | 2.8 |
| Cadmium | 6010B | ND | 0.57 |
| Chromium | 6010B | 32 | 0.57 |
| Lead | 6010B | ND | 5.7 |
| Mercury | 7471A | ND | 0.28 |
| Selenium | 6010B | ND | 11 |
| Silver | 6010B | ND | 0.57 |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A

Date Extracted: 3-13&14-08 Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-05 **Client ID: B-3_6**

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 6010B | ND | 11 |
| Barium | 6010B | 46 | 2.8 |
| Cadmium | 6010B | ND | 0.56 |
| Chromium | 6010B | 22 | 0.56 |
| Lead | 6010B | ND | 5.6 |
| Mercury | 7471A | ND | 0.28 |
| Selenium | 6010B | ND | 11 |
| Silver | 6010B | ND | 0.56 |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A METHOD BLANK QUALITY CONTROL

Date Extracted: 3-13&14-08
Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0313S2&MB0314S2

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 6010B | ND | 10 |
| Barium | 6010B | ND | 2.5 |
| Cadmium | 6010B | ND | 0.50 |
| Chromium | 6010B | ND | 0.50 |
| Lead | 6010B | ND | 5.0 |
| Mercury | 7471A | ND | 0.25 |
| Selenium | 6010B | ND | 10 |
| Silver | 6010B | ND | 0.50 |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A DUPLICATE QUALITY CONTROL

Date Extracted: 3-13&14-08 Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-087-19

| Analyte | Sample Result | Duplicate Result | RPD | PQL | Flags |
|----------|------------------|---------------------|-----|------|-------|
| Arsenic | 13.7 | 13.0 | 5 | 10 | |
| Barium | 87.8 | 84.6 | 4 | 2.5 | |
| Cadmium | ND | ND | NA | 0.50 | |
| Chromium | 35.0 | 33.0 | 6 | 0.50 | |
| Lead | 13.5 | 15.1 | 12 | 5.0 | |
| Mercury | ND | ND | NA | 0.25 | |
| Selenium | ND | ND | NA | 10 | |
| Silver | ND | ND | NA | 0.50 | |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A MS/MSD QUALITY CONTROL

Date Extracted: 3-13&14-08 Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-087-19

| Analyte | Spike Level | MS | Percent Recovery | MSD | Percent Recovery | RPD | Flags |
|----------|----------------|-------|---------------------|-------|---------------------|-----|-------|
| Arsenic | 100 | 103 | 90 | 105 | 91 | 2 | |
| Barium | 100 | 185 | 97 | 193 | 105 | 5 | |
| Cadmium | 50 | 47.6 | 95 | 48.2 | 96 | 1 | |
| Chromium | 100 | 129 | 94 | 128 | 93 | 0 | |
| Lead | 250 | 244 | 92 | 267 | 102 | 9 | |
| Mercury | 0.50 | 0.494 | 99 | 0.504 | 101 | 2 | |
| Selenium | 100 | 87.5 | 87 | 88.3 | 88 | 1 | |
| Silver | 25 | 21.9 | 88 | 22.1 | 88 | 1 | |

Project: 8915 162890

SOLUBLE HEXAVALENT CHROMIUM WATER EXTRACTION EPA 7196A

Project: 8915 162890

SOLUBLE HEXAVALENT CHROMIUM
WATER EXTRACTION
EPA 7196A
METHOD BLANK QUALITY CONTROL

Project: 8915 162890

SOLUBLE HEXAVALENT CHROMIUM
WATER EXTRACTION
EPA 7196A
DUPLICATE QUALITY CONTROL

Project: 8915 162890

SOLUBLE HEXAVALENT CHROMIUM
WATER EXTRACTION
EPA 7196A
MS/MSD QUALITY CONTROL

Project: 8915 162890

% MOISTURE

Date Analyzed: 3-12-08

| Client ID | Lab ID | % Moisture |
|-----------|-----------|------------|
| B-1_5 | 03-062-01 | 13 |
| B-1_3 | 03-062-02 | 12 |
| B-1_8 | 03-062-03 | 10 |
| B-3_3 | 03-062-04 | 19 |
| B-3_6 | 03-062-05 | 10 |



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-napthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical _____
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- Y Sample extract treated with an acid/silica gel cleanup procedure.
- Z Analyte may have been introduced by an outside source. See case narrative.
- ND Not Detected at PQL
- PQL Practical Quantitation Limit
- RPD Relative Percent Difference

Chain of Custody

% Moisture Afterlyze only if detections in Samples. HEN CHEWNE 0 Requested Analysis HEM py 1664 03-062 K **ICLP Metals** × (8) alsteM AROR listo? $\frac{\times}{\mathbb{X}}$ X Atčt8 yd sebicidreH X × Laboratory Number: MIS \ G07S8 vd sHA9 \times -£00√S8 yd selitslovime Islogenated Volatiles by 8260B Ø 89/5/ 8 3/10/08 × Volatiles by 8260B × × **XQ-H9TWN** Date × × NWTPH-Gx/BITEX имтрн-нспр (TPH analysis 5 working days) □ 3 Day 1 Day ス Standard (7 working days) Turnaround Request (in working days) (Check One) AMEC (other) 840 835 300 8/10/68 830 845 336 18 4S Same Day ☐ 2 Day Date Sampled Project Name No Trunsfor OnSite Environmental Inc. Project Manage: Lery Lyn Incuye Sampled by: FIELD BLANK -RIPBLANK Sample Identification Project Number: 8415/62890 AMEC Relinquished by Relinquished by Relinquished by Received by Company: Received by Received by

DISTRIBUTION LEGEND: White - OnSite Copy Yellow - Report Copy Pink - Client Copy

Reviewed by/Date

Reviewed by/Date

Chromatograms with final report



14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 26, 2008

Cherilyn Inouye AMEC Earth & Environmental, Inc. 5007 Pacific Highway East, Suite 5 Tacoma, WA 98424

Re: Analytical Data for Project 8-915-16341-0

Laboratory Reference No. 0803-105

Dear Cherilyn:

Enclosed are the analytical results and associated quality control data for samples submitted on March 14, 2008.

The standard policy of OnSite Environmental Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Project: 8-915-16341-0

Case Narrative

Samples were collected on March 14, 2008 and received by the laboratory on March 14, 2008. They were maintained at the laboratory at a temperature of 2°C to 6°C except as noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

Project: 8-915-16341-0

NWTPH-Gx

Date Extracted: 3-20-08 Date Analyzed: 3-20-08

Matrix: Water Units: ug/L (ppb)

 Client ID:
 MW-1
 MW-2

 Lab ID:
 03-105-01
 03-105-02

| | Result | Flags | PQL | Result | Flags | PQL |
|-----------------------------------|--------|-------|-----|--------|-------|-----|
| TPH-Gas | ND | | 100 | ND | | 100 |
| Surrogate Recovery: Fluorobenzene | 110% | | | 113% | | |

Project: 8-915-16341-0

NWTPH-Gx

Date Extracted: 3-20-08 Date Analyzed: 3-20-08

Matrix: Water Units: ug/L (ppb)

Client ID: MW-3 DUP-1
Lab ID: 03-105-03 03-105-04

| | Result | Flags | PQL | Result | Flags | PQL |
|-----------------------------------|--------|-------|-----|--------|-------|-----|
| TPH-Gas | ND | | 100 | ND | | 100 |
| Surrogate Recovery: Fluorobenzene | 112% | | | 112% | | |

Project: 8-915-16341-0

NWTPH-Gx

Date Extracted: 3-20-08 Date Analyzed: 3-20-08

Matrix: Water Units: ug/L (ppb)

Client ID: field blank_3
Lab ID: 03-105-05

Result Flags PQL

TPH-Gas ND 100

Surrogate Recovery:

Fluorobenzene 111%

Project: 8-915-16341-0

NWTPH-Gx METHOD BLANK QUALITY CONTROL

Date Extracted: 3-20-08 Date Analyzed: 3-20-08

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0320W3

Result Flags PQL

TPH-Gas ND 100

Surrogate Recovery:

Fluorobenzene 111%

Project: 8-915-16341-0

NWTPH-Gx DUPLICATE QUALITY CONTROL

Date Extracted: 3-20-08 Date Analyzed: 3-20-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-145-03 03-145-03

Original Duplicate RPD Flags

TPH-Gas ND ND NA

Surrogate Recovery:

Fluorobenzene 115% 114%

Project: 8-915-16341-0

NWTPH-Dx

Date Extracted: 3-24-08 Date Analyzed: 3-24-08

Matrix: Water Units: mg/L (ppm)

| Client ID: Lab ID: | MW-1 03-105-01 | MW-2 03-105-02 | MW-3 03-105-03 |
|-----------------------|--------------------------|--------------------------|--------------------------|
| Lab ID. | 03-103-01 | 03-103-02 | 03-103-03 |
| Diesel Range: | ND | ND | ND |
| PQL: | 0.27 | 0.26 | 0.27 |
| Identification: | | | |
| | | | |
| Lube Oil Range: | ND | ND | ND |
| PQL: | 0.43 | 0.42 | 0.43 |
| Identification: | | | |
| Surrogate Recovery | | | |
| o-Terphenyl: | 84% | 89% | 92% |
| Flags: | Υ | Υ | Υ |

Project: 8-915-16341-0

NWTPH-Dx

Date Extracted: 3-24-08 Date Analyzed: 3-24-08

Matrix: Water Units: mg/L (ppm)

 Client ID:
 DUP-1

 Lab ID:
 03-105-04

Diesel Range: **ND**PQL: 0.28

Identification: ---

Lube Oil Range: **ND**PQL: 0.45

Identification: ---

Surrogate Recovery

o-Terphenyl: 88%

Flags: Y

Project: 8-915-16341-0

NWTPH-Dx METHOD BLANK QUALITY CONTROL

Date Extracted: 3-24-08 Date Analyzed: 3-24-08

Matrix: Water Units: mg/L (ppm)

Lab ID: MB0324W1

Diesel Range: **ND**PQL: 0.25

Identification: ---

Lube Oil Range: **ND**PQL: 0.40

Identification: ---

Surrogate Recovery

o-Terphenyl: 84%

Flags: Y

Project: 8-915-16341-0

NWTPH-Dx DUPLICATE QUALITY CONTROL

Date Extracted: 3-24-08 Date Analyzed: 3-24-08

Matrix: Water Units: mg/L (ppm)

Lab ID: 03-105-01 03-105-01 DUP

Diesel Range: **ND ND** PQL: 0.27 0.26

RPD: N/A

Surrogate Recovery

o-Terphenyl: 84% 81%

Flags: Y Y

Project: 8-915-16341-0

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-19-08 Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| Dichlorodifluoromethane | ND | | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | ND | | 0.20 |
| Acetone | ND | | 5.0 |
| lodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | ND | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | ND | | 0.20 |
| Vinyl Acetate | ND | | 1.0 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | ND | | 0.20 |
| 1,1,1-Trichloroethane | ND | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | ND | | 0.20 |
| 1,2-Dichloroethane | ND | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B Page 2 of 2

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | _ | 0.20 |
| Tetrachloroethene | ND | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | 0.32 | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | 0.30 | | 0.20 |
| m,p-Xylene | 0.75 | | 0.40 |
| o-Xylene | 0.30 | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 86 | 71-126 |
| Toluene-d8 | 84 | 76-116 |
| 4-Bromofluorobenzene | 76 | 70-123 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-19-08 Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| Dichlorodifluoromethane | ND | | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | 31 | | 0.20 |
| Acetone | ND | | 5.0 |
| Iodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | ND | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | 30 | | 0.20 |
| Vinyl Acetate | ND | | 1.0 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | 0.66 | | 0.20 |
| 1,1,1-Trichloroethane | 30 | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | 0.32 | | 0.20 |
| 1,2-Dichloroethane | 3.4 | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B Page 2 of 2

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | J | 0.20 |
| Tetrachloroethene | ND | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | ND | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | ND | | 0.20 |
| m,p-Xylene | ND | | 0.40 |
| o-Xylene | ND | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 85 | 71-126 |
| Toluene-d8 | 84 | 76-116 |
| 4-Bromofluorobenzene | 76 | 70-123 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-19-08 Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| Dichlorodifluoromethane | ND | | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | ND | | 0.20 |
| Acetone | ND | | 5.0 |
| lodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | ND | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | ND | | 0.20 |
| Vinyl Acetate | ND | | 1.0 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | ND | | 0.20 |
| 1,1,1-Trichloroethane | ND | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | ND | | 0.20 |
| 1,2-Dichloroethane | ND | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B Page 2 of 2

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | _ | 0.20 |
| Tetrachloroethene | 0.81 | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | ND | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | ND | | 0.20 |
| m,p-Xylene | ND | | 0.40 |
| o-Xylene | ND | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 88 | 71-126 |
| Toluene-d8 | 84 | 76-116 |
| 4-Bromofluorobenzene | 74 | 70-123 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-19-08 Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-105-04 **Client ID: DUP-1**

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| Dichlorodifluoromethane | ND | _ | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | ND | | 0.20 |
| Acetone | ND | | 5.0 |
| lodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | ND | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | ND | | 0.20 |
| Vinyl Acetate | ND | | 1.0 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | ND | | 0.20 |
| 1,1,1-Trichloroethane | ND | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | ND | | 0.20 |
| 1,2-Dichloroethane | ND | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B Page 2 of 2

Lab ID: 03-105-04 Client ID: DUP-1

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | J | 0.20 |
| Tetrachloroethene | ND | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | 0.31 | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | 0.28 | | 0.20 |
| m,p-Xylene | 0.75 | | 0.40 |
| o-Xylene | 0.29 | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |
| | | | |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 89 | 71-126 |
| Toluene-d8 | 85 | 76-116 |
| 4-Bromofluorobenzene | 76 | 70-123 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-19-08 Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-105-05
Client ID: Field blank_3

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| Dichlorodifluoromethane | ND | | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | ND | | 0.20 |
| Acetone | ND | | 5.0 |
| lodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | ND | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | ND | | 0.20 |
| Vinyl Acetate | ND | | 1.0 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | 1.2 | | 0.20 |
| 1,1,1-Trichloroethane | ND | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | ND | | 0.20 |
| 1,2-Dichloroethane | ND | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B Page 2 of 2

Lab ID: 03-105-05 Client ID: Field blank_3

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | | 0.20 |
| Tetrachloroethene | ND | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | ND | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | ND | | 0.20 |
| m,p-Xylene | ND | | 0.40 |
| o-Xylene | ND | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 89 | 71-126 |
| Toluene-d8 | 84 | 76-116 |
| 4-Bromofluorobenzene | 76 | 70-123 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 1 of 2

Date Extracted: 3-19-08 Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0319W1

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| Dichlorodifluoromethane | ND | _ | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | ND | | 0.20 |
| Acetone | ND | | 5.0 |
| Iodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | ND | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | ND | | 0.20 |
| Vinyl Acetate | ND | | 1.0 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | ND | | 0.20 |
| 1,1,1-Trichloroethane | ND | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | ND | | 0.20 |
| 1,2-Dichloroethane | ND | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 2 of 2

Lab ID: MB0319W1

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | | 0.20 |
| Tetrachloroethene | ND | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | ND | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | ND | | 0.20 |
| m,p-Xylene | ND | | 0.40 |
| o-Xylene | ND | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 88 | 71-126 |
| Toluene-d8 | 83 | 76-116 |
| 4-Bromofluorobenzene | 76 | 70-123 |

Project: 8-915-16341-0

VOLATILES by EPA 8260B SB/SBD QUALITY CONTROL

Date Extracted: 3-19-08
Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: SB0319W1

| | Spike | | Percent | | Percent | Recovery | |
|--------------------|--------|------|----------|------|----------|----------|-------|
| Compound | Amount | SB | Recovery | SBD | Recovery | Limits | Flags |
| | | | | | | | |
| 1,1-Dichloroethene | 10.0 | 8.91 | 89 | 8.16 | 82 | 70-130 | |
| Benzene | 10.0 | 9.55 | 96 | 9.55 | 96 | 70-130 | |
| Trichloroethene | 10.0 | 8.45 | 85 | 8.29 | 83 | 70-116 | |
| Toluene | 10.0 | 9.80 | 98 | 9.62 | 96 | 76-119 | |
| Chlorobenzene | 10.0 | 9.12 | 91 | 9.16 | 92 | 77-112 | |
| | | | | | | | |

| | RPD | | |
|--------------------|-----|-------|-------|
| | RPD | Limit | Flags |
| 1,1-Dichloroethene | 9 | 20 | |
| Benzene | 0 | 16 | |
| Trichloroethene | 2 | 16 | |
| Toluene | 2 | 15 | |
| Chlorobenzene | 0 | 15 | |
| | | | |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 1 of 3

Date Extracted: 3-17-08
Date Analyzed: 3-18&20-08

Matrix: Water Units: ug/L (ppb)

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|------|
| N-Nitrosodimethylamine | ND | _ | 1.0 |
| Pyridine | ND | | 1.0 |
| Phenol | ND | | 1.0 |
| Aniline | ND | | 1.0 |
| bis(2-Chloroethyl)ether | ND | | 1.0 |
| 2-Chlorophenol | ND | | 1.0 |
| 1,3-Dichlorobenzene | ND | | 1.0 |
| 1,4-Dichlorobenzene | ND | | 1.0 |
| Benzyl alcohol | ND | | 1.0 |
| 1,2-Dichlorobenzene | ND | | 1.0 |
| 2-Methylphenol (o-Cresol) | ND | | 1.0 |
| bis(2-Chloroisopropyl)ether | ND | | 1.0 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 1.0 |
| N-Nitroso-di-n-propylamine | ND | | 1.0 |
| Hexachloroethane | ND | | 1.0 |
| Nitrobenzene | ND | | 1.0 |
| Isophorone | ND | | 1.0 |
| 2-Nitrophenol | ND | | 1.0 |
| 2,4-Dimethylphenol | ND | | 1.0 |
| bis(2-Chloroethoxy)methane | ND | | 1.0 |
| 2,4-Dichlorophenol | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 1.0 |
| Naphthalene | ND | | 0.10 |
| 4-Chloroaniline | ND | | 1.0 |
| Hexachlorobutadiene | ND | | 1.0 |
| 4-Chloro-3-methylphenol | ND | | 1.0 |
| 2-Methylnaphthalene | ND | | 0.10 |
| 1-Methylnaphthalene | ND | | 0.10 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|------|
| Hexachlorocyclopentadiene | ND | • | 1.0 |
| 2,4,6-Trichlorophenol | ND | | 1.0 |
| 2,3-Dichloroaniline | ND | | 1.0 |
| 2,4,5-Trichlorophenol | ND | | 1.0 |
| 2-Chloronaphthalene | ND | | 1.0 |
| 2-Nitroaniline | ND | | 1.0 |
| 1,4-Dinitrobenzene | ND | | 1.0 |
| Dimethylphthalate | ND | | 1.0 |
| 1,3-Dinitrobenzene | ND | | 1.0 |
| 2,6-Dinitrotoluene | ND | | 1.0 |
| 1,2-Dinitrobenzene | ND | | 1.0 |
| Acenaphthylene | ND | | 0.10 |
| 3-Nitroaniline | ND | | 1.0 |
| 2,4-Dinitrophenol | ND | | 5.2 |
| Acenaphthene | ND | | 0.10 |
| 4-Nitrophenol | ND | | 1.0 |
| 2,4-Dinitrotoluene | ND | | 1.0 |
| Dibenzofuran | ND | | 1.0 |
| 2,3,4,6-Tetrachlorophenol | ND | | 1.0 |
| 2,3,5,6-Tetrachlorophenol | ND | | 1.0 |
| Diethylphthalate | ND | | 1.0 |
| 4-Chlorophenyl-phenylether | ND | | 1.0 |
| 4-Nitroaniline | ND | | 1.0 |
| Fluorene | ND | | 0.10 |
| 4,6-Dinitro-2-methylphenol | ND | | 5.2 |
| N-Nitrosodiphenylamine | ND | | 1.0 |
| 1,2-Diphenylhydrazine | ND | | 1.0 |
| 4-Bromophenyl-phenylether | ND | | 1.0 |
| Hexachlorobenzene | ND | | 1.0 |
| Pentachlorophenol | ND | | 5.2 |
| Phenanthrene | ND | | 0.10 |
| Anthracene | ND | | 0.10 |
| Carbazole | ND | | 1.0 |
| Di-n-butylphthalate | 2.2 | | 1.0 |
| Fluoranthene | ND | | 0.10 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 3 of 3

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|-------|
| Benzidine | ND | | 10 |
| Pyrene | ND | | 0.10 |
| Butylbenzylphthalate | ND | | 1.0 |
| bis-2-Ethylhexyladipate | ND | | 1.0 |
| 3,3'-Dichlorobenzidine | ND | | 1.0 |
| Benzo[a]anthracene | ND | | 0.010 |
| Chrysene | ND | | 0.010 |
| bis(2-Ethylhexyl)phthalate | ND | | 1.0 |
| Di-n-octylphthalate | ND | | 1.0 |
| Benzo[b]fluoranthene | ND | | 0.010 |
| Benzo[k]fluoranthene | ND | | 0.010 |
| Benzo[a]pyrene | ND | | 0.010 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.010 |
| Dibenz[a,h]anthracene | ND | | 0.010 |
| Benzo[g,h,i]perylene | ND | | 0.010 |

| Surrogate : | Percent | Control |
|----------------------|----------|----------|
| | Recovery | Limits |
| 2-Fluorophenol | 54 | 10 - 90 |
| Phenol-d6 | 46 | 35 - 100 |
| Nitrobenzene-d5 | 86 | 30 - 100 |
| 2-Fluorobiphenyl | 93 | 39 - 100 |
| 2,4,6-Tribromophenol | 102 | 50 - 105 |
| Terphenyl-d14 | 109 | 49 - 115 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 1 of 3

Date Extracted: 3-17-08
Date Analyzed: 3-18&20-08

Matrix: Water Units: ug/L (ppb)

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|-------|
| N-Nitrosodimethylamine | ND | • | 0.98 |
| Pyridine | ND | | 0.98 |
| Phenol | ND | | 0.98 |
| Aniline | ND | | 0.98 |
| bis(2-Chloroethyl)ether | ND | | 0.98 |
| 2-Chlorophenol | ND | | 0.98 |
| 1,3-Dichlorobenzene | ND | | 0.98 |
| 1,4-Dichlorobenzene | ND | | 0.98 |
| Benzyl alcohol | ND | | 0.98 |
| 1,2-Dichlorobenzene | ND | | 0.98 |
| 2-Methylphenol (o-Cresol) | ND | | 0.98 |
| bis(2-Chloroisopropyl)ether | ND | | 0.98 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 0.98 |
| N-Nitroso-di-n-propylamine | ND | | 0.98 |
| Hexachloroethane | ND | | 0.98 |
| Nitrobenzene | ND | | 0.98 |
| Isophorone | ND | | 0.98 |
| 2-Nitrophenol | ND | | 0.98 |
| 2,4-Dimethylphenol | ND | | 0.98 |
| bis(2-Chloroethoxy)methane | ND | | 0.98 |
| 2,4-Dichlorophenol | ND | | 0.98 |
| 1,2,4-Trichlorobenzene | ND | | 0.98 |
| Naphthalene | ND | | 0.098 |
| 4-Chloroaniline | ND | | 0.98 |
| Hexachlorobutadiene | ND | | 0.98 |
| 4-Chloro-3-methylphenol | ND | | 0.98 |
| 2-Methylnaphthalene | ND | | 0.098 |
| 1-Methylnaphthalene | ND | | 0.098 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|-------|
| Hexachlorocyclopentadiene | ND | J | 0.98 |
| 2,4,6-Trichlorophenol | ND | | 0.98 |
| 2,3-Dichloroaniline | ND | | 0.98 |
| 2,4,5-Trichlorophenol | ND | | 0.98 |
| 2-Chloronaphthalene | ND | | 0.98 |
| 2-Nitroaniline | ND | | 0.98 |
| 1,4-Dinitrobenzene | ND | | 0.98 |
| Dimethylphthalate | ND | | 0.98 |
| 1,3-Dinitrobenzene | ND | | 0.98 |
| 2,6-Dinitrotoluene | ND | | 0.98 |
| 1,2-Dinitrobenzene | ND | | 0.98 |
| Acenaphthylene | ND | | 0.098 |
| 3-Nitroaniline | ND | | 0.98 |
| 2,4-Dinitrophenol | ND | | 4.9 |
| Acenaphthene | ND | | 0.098 |
| 4-Nitrophenol | ND | | 0.98 |
| 2,4-Dinitrotoluene | ND | | 0.98 |
| Dibenzofuran | ND | | 0.98 |
| 2,3,4,6-Tetrachlorophenol | ND | | 0.98 |
| 2,3,5,6-Tetrachlorophenol | ND | | 0.98 |
| Diethylphthalate | ND | | 0.98 |
| 4-Chlorophenyl-phenylether | ND | | 0.98 |
| 4-Nitroaniline | ND | | 0.98 |
| Fluorene | ND | | 0.098 |
| 4,6-Dinitro-2-methylphenol | ND | | 4.9 |
| N-Nitrosodiphenylamine | ND | | 0.98 |
| 1,2-Diphenylhydrazine | ND | | 0.98 |
| 4-Bromophenyl-phenylether | ND | | 0.98 |
| Hexachlorobenzene | ND | | 0.98 |
| Pentachlorophenol | ND | | 4.9 |
| Phenanthrene | ND | | 0.098 |
| Anthracene | ND | | 0.098 |
| Carbazole | ND | | 0.98 |
| Di-n-butylphthalate | ND | | 0.98 |
| Fluoranthene | ND | | 0.098 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 3 of 3

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Benzidine | ND | | 9.8 |
| Pyrene | 0.17 | | 0.098 |
| Butylbenzylphthalate | ND | | 0.98 |
| bis-2-Ethylhexyladipate | ND | | 0.98 |
| 3,3'-Dichlorobenzidine | ND | | 0.98 |
| Benzo[a]anthracene | ND | | 0.0098 |
| Chrysene | ND | | 0.0098 |
| bis(2-Ethylhexyl)phthalate | ND | | 0.98 |
| Di-n-octylphthalate | ND | | 0.98 |
| Benzo[b]fluoranthene | ND | | 0.0098 |
| Benzo[k]fluoranthene | ND | | 0.0098 |
| Benzo[a]pyrene | ND | | 0.0098 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.0098 |
| Dibenz[a,h]anthracene | ND | | 0.0098 |
| Benzo[g,h,i]perylene | ND | | 0.0098 |

| Surrogate : | Percent Recovery | Control Limits |
|----------------------|---------------------|-------------------|
| 2-Fluorophenol | 49 | 10 - 90 |
| Phenol-d6 | 45 | 35 - 100 |
| Nitrobenzene-d5 | 80 | 30 - 100 |
| 2-Fluorobiphenyl | 92 | 39 - 100 |
| 2,4,6-Tribromophenol | 103 | 50 - 105 |
| Terphenyl-d14 | 114 | 49 - 115 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 1 of 3

Date Extracted: 3-17-08
Date Analyzed: 3-18&20-08

Matrix: Water Units: ug/L (ppb)

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|------|
| N-Nitrosodimethylamine | ND | _ | 1.1 |
| Pyridine | ND | | 1.1 |
| Phenol | ND | | 1.1 |
| Aniline | ND | | 1.1 |
| bis(2-Chloroethyl)ether | ND | | 1.1 |
| 2-Chlorophenol | ND | | 1.1 |
| 1,3-Dichlorobenzene | ND | | 1.1 |
| 1,4-Dichlorobenzene | ND | | 1.1 |
| Benzyl alcohol | ND | | 1.1 |
| 1,2-Dichlorobenzene | ND | | 1.1 |
| 2-Methylphenol (o-Cresol) | ND | | 1.1 |
| bis(2-Chloroisopropyl)ether | ND | | 1.1 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 1.1 |
| N-Nitroso-di-n-propylamine | ND | | 1.1 |
| Hexachloroethane | ND | | 1.1 |
| Nitrobenzene | ND | | 1.1 |
| Isophorone | ND | | 1.1 |
| 2-Nitrophenol | ND | | 1.1 |
| 2,4-Dimethylphenol | ND | | 1.1 |
| bis(2-Chloroethoxy)methane | ND | | 1.1 |
| 2,4-Dichlorophenol | ND | | 1.1 |
| 1,2,4-Trichlorobenzene | ND | | 1.1 |
| Naphthalene | ND | | 0.11 |
| 4-Chloroaniline | ND | | 1.1 |
| Hexachlorobutadiene | ND | | 1.1 |
| 4-Chloro-3-methylphenol | ND | | 1.1 |
| 2-Methylnaphthalene | ND | | 0.11 |
| 1-Methylnaphthalene | ND | | 0.11 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|------|
| Hexachlorocyclopentadiene | ND | _ | 1.1 |
| 2,4,6-Trichlorophenol | ND | | 1.1 |
| 2,3-Dichloroaniline | ND | | 1.1 |
| 2,4,5-Trichlorophenol | ND | | 1.1 |
| 2-Chloronaphthalene | ND | | 1.1 |
| 2-Nitroaniline | ND | | 1.1 |
| 1,4-Dinitrobenzene | ND | | 1.1 |
| Dimethylphthalate | ND | | 1.1 |
| 1,3-Dinitrobenzene | ND | | 1.1 |
| 2,6-Dinitrotoluene | ND | | 1.1 |
| 1,2-Dinitrobenzene | ND | | 1.1 |
| Acenaphthylene | ND | | 0.11 |
| 3-Nitroaniline | ND | | 1.1 |
| 2,4-Dinitrophenol | ND | | 5.4 |
| Acenaphthene | ND | | 0.11 |
| 4-Nitrophenol | ND | | 1.1 |
| 2,4-Dinitrotoluene | ND | | 1.1 |
| Dibenzofuran | ND | | 1.1 |
| 2,3,4,6-Tetrachlorophenol | ND | | 1.1 |
| 2,3,5,6-Tetrachlorophenol | ND | | 1.1 |
| Diethylphthalate | ND | | 1.1 |
| 4-Chlorophenyl-phenylether | ND | | 1.1 |
| 4-Nitroaniline | ND | | 1.1 |
| Fluorene | ND | | 0.11 |
| 4,6-Dinitro-2-methylphenol | ND | | 5.4 |
| N-Nitrosodiphenylamine | ND | | 1.1 |
| 1,2-Diphenylhydrazine | ND | | 1.1 |
| 4-Bromophenyl-phenylether | ND | | 1.1 |
| Hexachlorobenzene | ND | | 1.1 |
| Pentachlorophenol | ND | | 5.4 |
| Phenanthrene | ND | | 0.11 |
| Anthracene | ND | | 0.11 |
| Carbazole | ND | | 1.1 |
| Di-n-butylphthalate | ND | | 1.1 |
| Fluoranthene | ND | | 0.11 |
| | | | |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 3 of 3

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|-------|
| Benzidine | ND | _ | 11 |
| Pyrene | ND | | 0.11 |
| Butylbenzylphthalate | ND | | 1.1 |
| bis-2-Ethylhexyladipate | ND | | 1.1 |
| 3,3'-Dichlorobenzidine | ND | | 1.1 |
| Benzo[a]anthracene | ND | | 0.011 |
| Chrysene | ND | | 0.011 |
| bis(2-Ethylhexyl)phthalate | ND | | 1.1 |
| Di-n-octylphthalate | ND | | 1.1 |
| Benzo[b]fluoranthene | ND | | 0.011 |
| Benzo[k]fluoranthene | ND | | 0.011 |
| Benzo[a]pyrene | ND | | 0.011 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.011 |
| Dibenz[a,h]anthracene | ND | | 0.011 |
| Benzo[g,h,i]perylene | ND | | 0.011 |

| Surrogate : | Percent | Control |
|----------------------|----------|----------|
| | Recovery | Limits |
| 2-Fluorophenol | 56 | 10 - 90 |
| Phenol-d6 | 46 | 35 - 100 |
| Nitrobenzene-d5 | 84 | 30 - 100 |
| 2-Fluorobiphenyl | 83 | 39 - 100 |
| 2,4,6-Tribromophenol | 99 | 50 - 105 |
| Terphenyl-d14 | 110 | 49 - 115 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 1 of 3

Date Extracted: 3-17-08
Date Analyzed: 3-18&20-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-105-04 **Client ID: DUP-1**

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|------|
| N-Nitrosodimethylamine | ND | _ | 1.1 |
| Pyridine | ND | | 1.1 |
| Phenol | ND | | 1.1 |
| Aniline | ND | | 1.1 |
| bis(2-Chloroethyl)ether | ND | | 1.1 |
| 2-Chlorophenol | ND | | 1.1 |
| 1,3-Dichlorobenzene | ND | | 1.1 |
| 1,4-Dichlorobenzene | ND | | 1.1 |
| Benzyl alcohol | ND | | 1.1 |
| 1,2-Dichlorobenzene | ND | | 1.1 |
| 2-Methylphenol (o-Cresol) | ND | | 1.1 |
| bis(2-Chloroisopropyl)ether | ND | | 1.1 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 1.1 |
| N-Nitroso-di-n-propylamine | ND | | 1.1 |
| Hexachloroethane | ND | | 1.1 |
| Nitrobenzene | ND | | 1.1 |
| Isophorone | ND | | 1.1 |
| 2-Nitrophenol | ND | | 1.1 |
| 2,4-Dimethylphenol | ND | | 1.1 |
| bis(2-Chloroethoxy)methane | ND | | 1.1 |
| 2,4-Dichlorophenol | ND | | 1.1 |
| 1,2,4-Trichlorobenzene | ND | | 1.1 |
| Naphthalene | ND | | 0.11 |
| 4-Chloroaniline | ND | | 1.1 |
| Hexachlorobutadiene | ND | | 1.1 |
| 4-Chloro-3-methylphenol | ND | | 1.1 |
| 2-Methylnaphthalene | ND | | 0.11 |
| 1-Methylnaphthalene | ND | | 0.11 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

 Lab ID:
 03-105-04

 Client ID:
 DUP-1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|------|
| Hexachlorocyclopentadiene | ND | _ | 1.1 |
| 2,4,6-Trichlorophenol | ND | | 1.1 |
| 2,3-Dichloroaniline | ND | | 1.1 |
| 2,4,5-Trichlorophenol | ND | | 1.1 |
| 2-Chloronaphthalene | ND | | 1.1 |
| 2-Nitroaniline | ND | | 1.1 |
| 1,4-Dinitrobenzene | ND | | 1.1 |
| Dimethylphthalate | ND | | 1.1 |
| 1,3-Dinitrobenzene | ND | | 1.1 |
| 2,6-Dinitrotoluene | ND | | 1.1 |
| 1,2-Dinitrobenzene | ND | | 1.1 |
| Acenaphthylene | ND | | 0.11 |
| 3-Nitroaniline | ND | | 1.1 |
| 2,4-Dinitrophenol | ND | | 5.4 |
| Acenaphthene | ND | | 0.11 |
| 4-Nitrophenol | ND | | 1.1 |
| 2,4-Dinitrotoluene | ND | | 1.1 |
| Dibenzofuran | ND | | 1.1 |
| 2,3,4,6-Tetrachlorophenol | ND | | 1.1 |
| 2,3,5,6-Tetrachlorophenol | ND | | 1.1 |
| Diethylphthalate | ND | | 1.1 |
| 4-Chlorophenyl-phenylether | ND | | 1.1 |
| 4-Nitroaniline | ND | | 1.1 |
| Fluorene | ND | | 0.11 |
| 4,6-Dinitro-2-methylphenol | ND | | 5.4 |
| N-Nitrosodiphenylamine | ND | | 1.1 |
| 1,2-Diphenylhydrazine | ND | | 1.1 |
| 4-Bromophenyl-phenylether | ND | | 1.1 |
| Hexachlorobenzene | ND | | 1.1 |
| Pentachlorophenol | ND | | 5.4 |
| Phenanthrene | ND | | 0.11 |
| Anthracene | ND | | 0.11 |
| Carbazole | ND | | 1.1 |
| Di-n-butylphthalate | 2.3 | | 1.1 |
| Fluoranthene | ND | | 0.11 |
| | | | |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM

page 3 of 3

 Lab ID:
 03-105-04

 Client ID:
 DUP-1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|-------|
| Benzidine | ND | | 11 |
| Pyrene | ND | | 0.11 |
| Butylbenzylphthalate | ND | | 1.1 |
| bis-2-Ethylhexyladipate | ND | | 1.1 |
| 3,3'-Dichlorobenzidine | ND | | 1.1 |
| Benzo[a]anthracene | ND | | 0.011 |
| Chrysene | ND | | 0.011 |
| bis(2-Ethylhexyl)phthalate | ND | | 1.1 |
| Di-n-octylphthalate | ND | | 1.1 |
| Benzo[b]fluoranthene | ND | | 0.011 |
| Benzo[k]fluoranthene | ND | | 0.011 |
| Benzo[a]pyrene | ND | | 0.011 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.011 |
| Dibenz[a,h]anthracene | ND | | 0.011 |
| Benzo[g,h,i]perylene | ND | | 0.011 |

| Surrogate : | Percent Recovery | Control Limits |
|----------------------|---------------------|-------------------|
| 2-Fluorophenol | 52 | 10 - 90 |
| Phenol-d6 | 45 | 35 - 100 |
| Nitrobenzene-d5 | 78 | 30 - 100 |
| 2-Fluorobiphenyl | 87 | 39 - 100 |
| 2,4,6-Tribromophenol | 104 | 50 - 105 |
| Terphenyl-d14 | 113 | 49 - 115 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 1 of 3

Date Extracted: 3-17-08
Date Analyzed: 3-18&20-08

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0317W1

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|------|
| N-Nitrosodimethylamine | ND | 9- | 1.0 |
| Pyridine | ND | | 1.0 |
| Phenol | ND | | 1.0 |
| Aniline | ND | | 1.0 |
| bis(2-Chloroethyl)ether | ND | | 1.0 |
| 2-Chlorophenol | ND | | 1.0 |
| 1,3-Dichlorobenzene | ND | | 1.0 |
| 1,4-Dichlorobenzene | ND | | 1.0 |
| Benzyl alcohol | ND | | 1.0 |
| 1,2-Dichlorobenzene | ND | | 1.0 |
| 2-Methylphenol (o-Cresol) | ND | | 1.0 |
| bis(2-Chloroisopropyl)ether | ND | | 1.0 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 1.0 |
| N-Nitroso-di-n-propylamine | ND | | 1.0 |
| Hexachloroethane | ND | | 1.0 |
| Nitrobenzene | ND | | 1.0 |
| Isophorone | ND | | 1.0 |
| 2-Nitrophenol | ND | | 1.0 |
| 2,4-Dimethylphenol | ND | | 1.0 |
| bis(2-Chloroethoxy)methane | ND | | 1.0 |
| 2,4-Dichlorophenol | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 1.0 |
| Naphthalene | ND | | 0.10 |
| 4-Chloroaniline | ND | | 1.0 |
| Hexachlorobutadiene | ND | | 1.0 |
| 4-Chloro-3-methylphenol | ND | | 1.0 |
| 2-Methylnaphthalene | ND | | 0.10 |
| 1-Methylnaphthalene | ND | | 0.10 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 2 of 3

Lab ID: MB0317W1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|------|
| Hexachlorocyclopentadiene | ND | | 1.0 |
| 2,4,6-Trichlorophenol | ND | | 1.0 |
| 2,3-Dichloroaniline | ND | | 1.0 |
| 2,4,5-Trichlorophenol | ND | | 1.0 |
| 2-Chloronaphthalene | ND | | 1.0 |
| 2-Nitroaniline | ND | | 1.0 |
| 1,4-Dinitrobenzene | ND | | 1.0 |
| Dimethylphthalate | ND | | 1.0 |
| 1,3-Dinitrobenzene | ND | | 1.0 |
| 2,6-Dinitrotoluene | ND | | 1.0 |
| 1,2-Dinitrobenzene | ND | | 1.0 |
| Acenaphthylene | ND | | 0.10 |
| 3-Nitroaniline | ND | | 1.0 |
| 2,4-Dinitrophenol | ND | | 5.0 |
| Acenaphthene | ND | | 0.10 |
| 4-Nitrophenol | ND | | 1.0 |
| 2,4-Dinitrotoluene | ND | | 1.0 |
| Dibenzofuran | ND | | 1.0 |
| 2,3,4,6-Tetrachlorophenol | ND | | 1.0 |
| 2,3,5,6-Tetrachlorophenol | ND | | 1.0 |
| Diethylphthalate | ND | | 1.0 |
| 4-Chlorophenyl-phenylether | ND | | 1.0 |
| 4-Nitroaniline | ND | | 1.0 |
| Fluorene | ND | | 0.10 |
| 4,6-Dinitro-2-methylphenol | ND | | 5.0 |
| N-Nitrosodiphenylamine | ND | | 1.0 |
| 1,2-Diphenylhydrazine | ND | | 1.0 |
| 4-Bromophenyl-phenylether | ND | | 1.0 |
| Hexachlorobenzene | ND | | 1.0 |
| Pentachlorophenol | ND | | 5.0 |
| Phenanthrene | ND | | 0.10 |
| Anthracene | ND | | 0.10 |
| Carbazole | ND | | 1.0 |
| Di-n-butylphthalate | ND | | 1.0 |
| Fluoranthene | ND | | 0.10 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 3 of 3

Lab ID: MB0317W1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|-------|
| Benzidine | ND | | 10 |
| Pyrene | ND | | 0.10 |
| Butylbenzylphthalate | ND | | 1.0 |
| bis-2-Ethylhexyladipate | ND | | 1.0 |
| 3,3'-Dichlorobenzidine | ND | | 1.0 |
| Benzo[a]anthracene | ND | | 0.010 |
| Chrysene | ND | | 0.010 |
| bis(2-Ethylhexyl)phthalate | ND | | 1.0 |
| Di-n-octylphthalate | ND | | 1.0 |
| Benzo[b]fluoranthene | ND | | 0.010 |
| Benzo[k]fluoranthene | ND | | 0.010 |
| Benzo[a]pyrene | ND | | 0.010 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.010 |
| Dibenz[a,h]anthracene | ND | | 0.010 |
| Benzo[g,h,i]perylene | ND | | 0.010 |

| Surrogate : | Percent | Control |
|----------------------|----------|----------|
| _ | Recovery | Limits |
| 2-Fluorophenol | 54 | 10 - 90 |
| Phenol-d6 | 45 | 35 - 100 |
| Nitrobenzene-d5 | 84 | 30 - 100 |
| 2-Fluorobiphenyl | 86 | 39 - 100 |
| 2,4,6-Tribromophenol | 92 | 50 - 105 |
| Terphenyl-d14 | 110 | 49 - 115 |

Project: 8-915-16341-0

SEMIVOLATILES by EPA 8270D/SIM SB/SBD QUALITY CONTROL

Date Extracted: 3-17-08
Date Analyzed: 3-18-08

Matrix: Water Units: ug/L (ppb)

Lab ID: SB0317W1

| | Spike | | Percent | | Percent I | Recovery | |
|----------------------------|--------|------|----------|------|-----------|----------|-------|
| Compound: | Amount | SB | Recovery | SBD | Recovery | Limits | Flags |
| Phenol | 40.0 | 18.3 | 46 | 18.9 | 47 | 21-75 | |
| 2-Chlorophenol | 40.0 | 31.0 | 77 | 33.3 | 83 | 43-96 | |
| 1,4-Dichlorobenzene | 20.0 | 11.7 | 59 | 13.3 | 66 | 38-80 | |
| N-Nitroso-di-n-propylamine | 20.0 | 16.8 | 84 | 16.7 | 83 | 36-99 | |
| 1,2,4-Trichlorobenzene | 20.0 | 11.7 | 58 | 12.2 | 61 | 39-85 | |
| 4-Chloro-3-methylphenol | 40.0 | 36.8 | 92 | 36.2 | 91 | 50-105 | |
| Acenaphthene | 20.0 | 16.7 | 83 | 16.4 | 82 | 46-90 | |
| 2,4-Dinitrotoluene | 20.0 | 20.1 | 100 | 19.5 | 98 | 50-122 | |
| 4-Nitrophenol | 40.0 | 26.6 | 66 | 25.4 | 64 | 30-116 | |
| Pentachlorophenol | 40.0 | 36.7 | 92 | 36.4 | 91 | 40-112 | |
| Pyrene | 20.0 | 20.5 | 103 | 20.1 | 101 | 51-105 | |

| | RPD | RPD Limits | Flags |
|----------------------------|-----|---------------|-------|
| Phenol | 3 | 31 | |
| 2-Chlorophenol | 7 | 29 | |
| 1,4-Dichlorobenzene | 13 | 33 | |
| N-Nitroso-di-n-propylamine | 1 | 28 | |
| 1,2,4-Trichlorobenzene | 4 | 32 | |
| 4-Chloro-3-methylphenol | 2 | 23 | |
| Acenaphthene | 2 | 24 | |
| 2,4-Dinitrotoluene | 3 | 30 | |
| 4-Nitrophenol | 4 | 30 | |
| Pentachlorophenol | 1 | 30 | |
| Pyrene | 2 | 18 | |

Project: 8-915-16341-0

PCBs by EPA 8082

Matrix: Water
Units: ug/L (ppb)

DCB

| | | | | Date | Date | |
|----------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | MW-1 | | | | | |
| Laboratory ID: | 03-105-01 | | | | | |
| Aroclor 1016 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1221 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1232 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1242 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1248 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1254 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1260 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1262 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1268 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| DCB | 77 | 45-121 | | | | |
| Client ID: | MW-2 | | | | | |
| Laboratory ID: | 03-105-02 | | | | | |
| Aroclor 1016 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1221 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1232 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1242 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1248 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1254 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1260 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1262 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1268 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| DCB | 99 | 45-121 | | | | |
| Client ID: | MW-3 | | | | | |
| _aboratory ID: | 03-105-03 | | | | | |
| Aroclor 1016 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | , |
| Aroclor 1221 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1232 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1242 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1248 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1254 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1260 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1262 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1268 | ND | 0.048 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| | | | | | | |

45-121

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Project: 8-915-16341-0

PCBs by EPA 8082

Matrix: Water
Units: ug/L (ppb)

| | | | | Date | Date | |
|----------------|-----------|-------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | DUP-1 | | | | | |
| Laboratory ID: | 03-105-04 | | | | | |
| Aroclor 1016 | ND | 0.058 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1221 | ND | 0.058 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1232 | ND | 0.058 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1242 | ND | 0.058 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1248 | ND | 0.058 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1254 | ND | 0.058 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1260 | ND | 0.058 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1262 | ND | 0.058 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1268 | ND | 0.058 | EPA 8082 | 3-18-08 | 3-18-08 | |

Surrogate: Percent Recovery Control Limits DCB 89 45-121

Project: 8-915-16341-0

PCBs by EPA 8082 QUALITY CONTROL

Matrix: Water
Units: ug/L (ppb)

| | | | | Date | Date | |
|----------------|---------------------------------------|-------|---------------------------------------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| METHOD BLANK | | | | | | |
| Laboratory ID: | MB0318W1 | | | | | |
| Aroclor 1016 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1221 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1232 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1242 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1248 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1254 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1260 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1262 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| Aroclor 1268 | ND | 0.050 | EPA 8082 | 3-18-08 | 3-18-08 | |
| • | · · · · · · · · · · · · · · · · · · · | | · · · · · · · · · · · · · · · · · · · | | | |

Surrogate: Percent Recovery Control Limits DCB 90 45-121

| Analyte | Re | sult | Spike | Level | Source Result | | cent overy | Recovery Limits | RPD | RPD Limit | Flags |
|----------------|-------|-------|-------|-------|------------------|-----|---------------|--------------------|-----|--------------|-------|
| SPIKE BLANKS | | | | | | | | | | | |
| Laboratory ID: | SB03 | 318W1 | | | | | | | | | |
| | SB | SBD | SB | SBD | | SB | SBD | | | | |
| Aroclor 1260 | 0.471 | 0.471 | 0.500 | 0.500 | N/A | 94 | 94 | 57-115 | 0 | 12 | _ |
| Surrogate: | | | | | | | | | | | |
| DCB | | | | | | 100 | 97 | 45-121 | | | |

Project: 8-915-16341-0

ORGANOCHLORINE PESTICIDES by EPA 8081A

Matrix: Water
Units: ug/L (ppb)

| 3 (11) | | | | Date | Date | |
|--------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | MW-1 | | | | | |
| Laboratory ID: | 03-105-01 | | | | | |
| alpha-BHC | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-BHC | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| beta-BHC | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| delta-BHC | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Aldrin | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor Epoxide | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-Chlordane | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| alpha-Chlordane | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDE | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan I | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Dieldrin | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDD | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan II | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDT | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Aldehyde | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Methoxychlor | ND | 0.011 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endsulfan Sulfate | ND | 0.0056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Ketone | ND | 0.023 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Toxaphene | ND | 0.056 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Surrogate: | Percent Recovery | Control Limits | _ | | _ | |
| | | | | | | |

Surrogate: Percent Recovery Control Limit
TCMX 61 30-105
DCB 80 30-121

Project: 8-915-16341-0

ORGANOCHLORINE PESTICIDES by EPA 8081A

Matrix: Water
Units: ug/L (ppb)

| 3 (11) | | | | Date | Date | |
|--------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | MW-2 | | | | | |
| Laboratory ID: | 03-105-02 | | | | | |
| alpha-BHC | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-BHC | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| beta-BHC | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| delta-BHC | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Aldrin | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor Epoxide | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-Chlordane | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| alpha-Chlordane | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDE | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan I | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Dieldrin | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDD | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan II | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDT | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Aldehyde | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Methoxychlor | ND | 0.0099 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endsulfan Sulfate | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Ketone | ND | 0.020 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Toxaphene | ND | 0.050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| TO10/ | - | 00.405 | | | | |

Surrogate: Percent Recovery Control Limit
TCMX 66 30-105
DCB 77 30-121

Project: 8-915-16341-0

ORGANOCHLORINE PESTICIDES by EPA 8081A

Matrix: Water Units: ug/L (ppb)

| 3 (11 / | | | | Date | Date | |
|--------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | MW-3 | | | | | |
| Laboratory ID: | 03-105-03 | | | | | |
| alpha-BHC | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-BHC | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| oeta-BHC | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| delta-BHC | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Aldrin | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor Epoxide | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-Chlordane | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| alpha-Chlordane | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 1,4'-DDE | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan I | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Dieldrin | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 1,4'-DDD | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan II | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 1,4'-DDT | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Aldehyde | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Methoxychlor | ND | 0.010 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endsulfan Sulfate | ND | 0.0051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Ketone | ND | 0.020 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Гохарhene | ND | 0.051 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| | | | | | | |

Surrogate:Percent RecoveryControl LimitTCMX6130-105DCB7330-121

Project: 8-915-16341-0

ORGANOCHLORINE PESTICIDES by EPA 8081A

Matrix: Water Units: ug/L (ppb)

| Analyto | Result | PQL | Method | Date Prepared | Date Analyzed | Elage |
|-----------------------|------------------|----------------|----------|------------------|------------------|-------|
| Analyte Client ID: | DUP-1 | FQL | Metriou | Frepareu | Allalyzeu | Flags |
| Laboratory ID: | 03-105-04 | | | | | |
| alpha-BHC | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-BHC | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| beta-BHC | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| delta-BHC | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Aldrin | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor Epoxide | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-Chlordane | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| alpha-Chlordane | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDE | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan I | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Dieldrin | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDD | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan II | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDT | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Aldehyde | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Methoxychlor | ND | 0.0096 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endsulfan Sulfate | ND | 0.0048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Ketone | ND | 0.019 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Toxaphene | ND | 0.048 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |

Surrogate:Percent RecoveryControl LimitTCMX6130-105DCB8530-121

Project: 8-915-16341-0

ORGANOCHLORINE PESTICIDES by EPA 8081A **QUALITY CONTROL**

Matrix: Water Units: ug/L (ppb)

| 3 (11) | | | | Date | Date | |
|--------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| METHOD BLANK | | | | | | |
| Laboratory ID: | MB0318W1 | | | | | |
| alpha-BHC | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-BHC | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| beta-BHC | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| delta-BHC | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Aldrin | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Heptachlor Epoxide | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| gamma-Chlordane | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| alpha-Chlordane | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDE | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan I | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Dieldrin | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDD | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endosulfan II | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| 4,4'-DDT | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Aldehyde | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Methoxychlor | ND | 0.010 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endsulfan Sulfate | ND | 0.0050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Endrin Ketone | ND | 0.020 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Toxaphene | ND | 0.050 | EPA 8081 | 3-18-08 | 3-19-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| TCMX | 67 | 30-105 | | | | |
| DOD | 0.7 | 20.424 | | | | |

| Surrogate: | Percent Recovery | Control Limits |
|------------|------------------|----------------|
| TCMX | 67 | 30-105 |
| DCB | 87 | 30-121 |

| Analyte | Res | sult | Snike | Level | Source Result | | rcent | Recovery Limits | RPD | RPD Limit | Flags |
|----------------|--------|----------|--------|--------|------------------|----|-------------|--------------------|-----|--------------|-------|
| SPIKE BLANKS | | <u> </u> | Орико | | rtoount | | , o v o . y | | | | age |
| Laboratory ID: | SB03 | 18W1 | | | | | | | | | |
| | SB | SBD | SB | SBD | | SB | SBD | | | | |
| gamma-BHC | 0.0427 | 0.0421 | 0.0500 | 0.0500 | N/A | 85 | 84 | 50-104 | 1 | 20 | |
| Heptachlor | 0.0360 | 0.0365 | 0.0500 | 0.0500 | N/A | 72 | 73 | 41-102 | 1 | 22 | |
| Aldrin | 0.0355 | 0.0351 | 0.0500 | 0.0500 | N/A | 71 | 70 | 29-101 | 1 | 25 | |
| Dieldrin | 0.101 | 0.100 | 0.125 | 0.125 | N/A | 81 | 80 | 55-110 | 1 | 24 | |
| Endrin | 0.109 | 0.109 | 0.125 | 0.125 | N/A | 87 | 87 | 54-120 | 0 | 22 | |
| 4,4'-DDT | 0.104 | 0.106 | 0.125 | 0.125 | N/A | 83 | 85 | 56-119 | 2 | 26 | |
| Surrogate: | | | | | | | | | | | |
| TCMX | | | | | | 67 | 66 | 30-105 | | | |
| DCB | | | | | | 79 | 79 | 30-121 | | | |

OnSite Environmental, Inc. 14648 NE 95th Street, Redmond, WA 98052 (425) 883-3881

Project: 8-915-16341-0

CHLORINATED ACID HERBICIDES by EPA 8151A

Matrix: Water
Units: ug/L (ppb)

2,4,5-TP (Silvex)

2,4,5-T

2,4-DB

Dinoseb

| 3 (11 -) | | | | Date | Date | |
|-------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | MW-1 | | | | | |
| Laboratory ID: | 03-105-01 | | | | | |
| Dalapon | ND | 0.23 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 4.6 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 4.6 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 0.46 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 0.0094 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-T | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-DB | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dinoseb | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| DCAA | 56 | 44-116 | | | | |
| | | | | | | |
| Client ID: | MW-2 | | | | | |
| Laboratory ID: | 03-105-02 | | | | | |
| Dalapon | ND | 0.22 | EPA 8151 | 3-21-08 | 3-24-08 | |
| Dicamba | ND | 0.023 | EPA 8151 | 3-21-08 | 3-24-08 | |
| MCPP | ND | 4.6 | EPA 8151 | 3-21-08 | 3-24-08 | |
| MCPA | ND | 4.6 | EPA 8151 | 3-21-08 | 3-24-08 | |
| Dichlorprop | ND | 0.023 | EPA 8151 | 3-21-08 | 3-24-08 | |
| 2,4-D | ND | 0.46 | EPA 8151 | 3-21-08 | 3-24-08 | |
| Pentachlorophenol | ND | 0.0093 | EPA 8151 | 3-21-08 | 3-24-08 | |

Surrogate: Percent Recovery Control Limits DCAA 76 44-116

ND

ND

ND

ND

0.023

0.023

0.023

0.023

EPA 8151

EPA 8151

EPA 8151

EPA 8151

3-21-08

3-21-08

3-21-08

3-21-08

3-24-08

3-24-08

3-24-08

3-24-08

Project: 8-915-16341-0

CHLORINATED ACID HERBICIDES by EPA 8151A

Matrix: Water
Units: ug/L (ppb)

2,4,5-T

2,4-DB

Dinoseb

| - 3 - (i.i7) | | | | Date | Date | |
|---------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | MW-3 | | | | | |
| Laboratory ID: | 03-105-03 | | | | | |
| Dalapon | ND | 0.22 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 4.5 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 4.5 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 0.45 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 0.0091 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-T | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-DB | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dinoseb | ND | 0.023 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| DCAA | 67 | 44-116 | | | | |
| Client ID: | DUP-1 | | | | | |
| Laboratory ID: | 03-105-04 | | | | | |
| Dalapon | ND | 0.23 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 0.024 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 4.7 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 4.7 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 0.024 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 0.47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 0.0095 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 0.024 | EPA 8151 | 3-18-08 | 3-18-08 | |
| ' | | | | | | |

Surrogate: Percent Recovery Control Limits DCAA 76 44-116

ND

ND

ND

0.024

0.024

0.024

EPA 8151

EPA 8151

EPA 8151

3-18-08

3-18-08

3-18-08

3-18-08

3-18-08

3-18-08

Project: 8-915-16341-0

CHLORINATED ACID HERBICIDES by EPA 8151A METHOD BLANK QUALITY CONTROL

Matrix: Water
Units: ug/L (ppb)

| | | | | Date | Date | |
|-------------------|----------|--------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Laboratory ID: | MB0318W1 | | | | | |
| Dalapon | ND | 0.23 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 0.024 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 4.7 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 4.7 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 0.024 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 0.47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 0.0095 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 0.024 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-T | ND | 0.024 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-DB | ND | 0.024 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dinoseb | ND | 0.024 | EPA 8151 | 3-18-08 | 3-18-08 | |

Surrogate: Percent Recovery Control Limits DCAA 74 44-116

| Laboratory ID: | MB0321W1 | | | | | |
|-------------------|----------|--------|----------|---------|---------|---|
| Dalapon | ND | 0.23 | EPA 8151 | 3-21-08 | 3-24-08 | _ |
| Dicamba | ND | 0.024 | EPA 8151 | 3-21-08 | 3-24-08 | |
| MCPP | ND | 4.7 | EPA 8151 | 3-21-08 | 3-24-08 | |
| MCPA | ND | 4.7 | EPA 8151 | 3-21-08 | 3-24-08 | |
| Dichlorprop | ND | 0.024 | EPA 8151 | 3-21-08 | 3-24-08 | |
| 2,4-D | ND | 0.47 | EPA 8151 | 3-21-08 | 3-24-08 | |
| Pentachlorophenol | ND | 0.0095 | EPA 8151 | 3-21-08 | 3-24-08 | |
| 2,4,5-TP (Silvex) | ND | 0.024 | EPA 8151 | 3-21-08 | 3-24-08 | |
| 2,4,5-T | ND | 0.024 | EPA 8151 | 3-21-08 | 3-24-08 | |
| 2,4-DB | ND | 0.024 | EPA 8151 | 3-21-08 | 3-24-08 | |
| Dinoseb | ND | 0.024 | EPA 8151 | 3-21-08 | 3-24-08 | |

Surrogate: Percent Recovery Control Limits DCAA 72 44-116

Project: 8-915-16341-0

CHLORINATED ACID HERBICIDES by EPA 8151A SB/SBD QUALITY CONTROL

Matrix: Water
Units: ug/L (ppb)

| | | | | | Source | Per | cent | Recovery | | RPD | |
|----------------|--------|-------|-------|-------------|--------|----------|------|----------|-----|-------|-------|
| Analyte | Result | | Spike | Spike Level | | Recovery | | Limits | RPD | Limit | Flags |
| Laboratory ID: | SB03 | 18W1 | | | | | | | | | |
| | SB | SBD | SB | SBD | | SB | SBD | | | | |
| Dicamba | 0.824 | 0.797 | 1.00 | 1.00 | N/A | 82 | 80 | 29-121 | 3 | 20 | |
| 2,4-D | 0.797 | 0.786 | 1.00 | 1.00 | N/A | 80 | 79 | 47-110 | 1 | 20 | |
| 2,4,5-T | 0.897 | 0.868 | 1.00 | 1.00 | N/A | 90 | 87 | 61-119 | 3 | 20 | |
| 2,4-DB | 0.942 | 1.11 | 1.00 | 1.00 | N/A | 94 | 111 | 70-144 | 16 | 20 | |
| Surrogate: | | | | | | | | | | | |
| DCAA | | | | | | 78 | 79 | 44-116 | | | |
| | | | | | | | | | | | |
| Laboratory ID: | SB03 | 21W1 | | | | | | | | | |
| | SB | SBD | SB | SBD | | SB | SBD | | | | |
| Dicamba | 0.615 | 0.654 | 1.00 | 1.00 | N/A | 61 | 65 | 29-121 | 6 | 20 | |
| 2,4-D | 0.806 | 0.787 | 1.00 | 1.00 | N/A | 81 | 79 | 47-110 | 2 | 20 | |
| 2,4,5-T | 0.912 | 0.902 | 1.00 | 1.00 | N/A | 91 | 90 | 61-119 | 1 | 20 | |
| 2,4-DB | 0.971 | 0.950 | 1.00 | 1.00 | N/A | 97 | 95 | 70-144 | 2 | 20 | |
| Surrogate: | | | | | | | | | | | |
| DCAA | | | | | | 75 | 78 | 44-116 | | | |

Project: 8-915-16341-0

DISSOLVED METALS EPA 200.8/7470A

Date Filtered: 3-17-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-105-01 Client ID: MW-1

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 200.8 | ND | 3.0 |
| Barium | 200.8 | 33 | 25 |
| Cadmium | 200.8 | ND | 4.0 |
| Chromium | 200.8 | ND | 10 |
| Lead | 200.8 | ND | 1.0 |
| Mercury | 7470A | ND | 0.50 |
| Selenium | 200.8 | ND | 5.0 |
| Silver | 200.8 | ND | 10 |

Project: 8-915-16341-0

DISSOLVED METALS EPA 200.8/7470A

Date Filtered: 3-17-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-105-02 Client ID: MW-2

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 200.8 | ND | 3.0 |
| Barium | 200.8 | ND | 25 |
| Cadmium | 200.8 | ND | 4.0 |
| Chromium | 200.8 | ND | 10 |
| Lead | 200.8 | ND | 1.0 |
| Mercury | 7470A | ND | 0.50 |
| Selenium | 200.8 | ND | 5.0 |
| Silver | 200.8 | ND | 10 |

Project: 8-915-16341-0

DISSOLVED METALS EPA 200.8/7470A

Date Filtered: 3-17-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-105-03 Client ID: MW-3

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 200.8 | ND | 3.0 |
| Barium | 200.8 | 36 | 25 |
| Cadmium | 200.8 | ND | 4.0 |
| Chromium | 200.8 | ND | 10 |
| Lead | 200.8 | ND | 1.0 |
| Mercury | 7470A | ND | 0.50 |
| Selenium | 200.8 | ND | 5.0 |
| Silver | 200.8 | ND | 10 |

Project: 8-915-16341-0

DISSOLVED METALS EPA 200.8/7470A

Date Filtered: 3-17-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-105-04 Client ID: DUP-1

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 200.8 | ND | 3.0 |
| Barium | 200.8 | 31 | 25 |
| Cadmium | 200.8 | ND | 4.0 |
| Chromium | 200.8 | ND | 10 |
| Lead | 200.8 | ND | 1.0 |
| Mercury | 7470A | ND | 0.50 |
| Selenium | 200.8 | ND | 5.0 |
| Silver | 200.8 | ND | 10 |

Project: 8-915-16341-0

DISSOLVED METALS EPA 200.8/7470A METHOD BLANK QUALITY CONTROL

Date Filtered: 3-17-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0317D1

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 200.8 | ND | 3.0 |
| Barium | 200.8 | ND | 25 |
| Cadmium | 200.8 | ND | 4.0 |
| Chromium | 200.8 | ND | 10 |
| Lead | 200.8 | ND | 1.0 |
| Mercury | 7470A | ND | 0.50 |
| Selenium | 200.8 | ND | 5.0 |
| Silver | 200.8 | ND | 10 |

Project: 8-915-16341-0

DISSOLVED METALS EPA 200.8/7470A DUPLICATE QUALITY CONTROL

Date Filtered: 3-17-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-077-01

| Analyte | Sample Result | Duplicate Result | RPD | PQL | Flags |
|----------|------------------|---------------------|-----|------|-------|
| Arsenic | ND | ND | NA | 3.0 | |
| Barium | 33.5 | 32.1 | 4 | 25 | |
| Cadmium | ND | ND | NA | 4.0 | |
| Chromium | ND | ND | NA | 10 | |
| Lead | ND | ND | NA | 1.0 | |
| Mercury | ND | ND | NA | 0.50 | |
| Selenium | ND | ND | NA | 5.0 | |
| Silver | ND | ND | NA | 10 | |

Project: 8-915-16341-0

DISSOLVED METALS EPA 200.8/7470A MS/MSD QUALITY CONTROL

Date Filtered: 3-17-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-077-01

| Analyte | Spike Level | MS | Percent Recovery | MSD | Percent Recovery | RPD | Flags |
|----------|----------------|------|---------------------|------|---------------------|-----|-------|
| Arsenic | 200 | 217 | 109 | 213 | 107 | 2 | |
| Barium | 200 | 231 | 99 | 234 | 100 | 1 | |
| Cadmium | 200 | 204 | 102 | 205 | 103 | 1 | |
| Chromium | 200 | 194 | 97 | 199 | 100 | 3 | |
| Lead | 200 | 197 | 99 | 196 | 98 | 1 | |
| Mercury | 12.5 | 12.4 | 99 | 12.4 | 99 | 0 | |
| Selenium | 200 | 223 | 112 | 219 | 110 | 2 | |
| Silver | 200 | 170 | 85 | 173 | 87 | 2 | |



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-napthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical ______
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- Y Sample extract treated with an acid/silica gel cleanup procedure.

Z -

- ND Not Detected at PQL
- PQL Practical Quantitation Limit
- RPD Relative Percent Difference

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| AMEC FARNCHENICONMENTAL Project Number: | W | |
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| 2 MW-Z | | |
| 3 MW-3 | 1 020 | |
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14648 NE 95th Street, Redmond, WA 98052 • (425) 883-3881

March 25, 2008

Cherilyn Inouye AMEC Earth & Environmental, Inc. 5007 Pacific Highway East, Suite 5 Tacoma, WA 98424

Re: Analytical Data for Project 8915 162890

Laboratory Reference No. 0803-076

Dear Cherilyn:

Enclosed are the analytical results and associated quality control data for samples submitted on March 12, 2008.

The standard policy of OnSite Environmental Inc. is to store your samples for 30 days from the date of receipt. If you require longer storage, please contact the laboratory.

We appreciate the opportunity to be of service to you on this project. If you have any questions concerning the data, or need additional information, please feel free to call me.

Sincerely,

David Baumeister Project Manager

Enclosures

Project: 8915 162890

Case Narrative Page 1 of 2

Samples were collected on March 11, 2008 and received by the laboratory on March 12, 2008. They were maintained at the laboratory at a temperature of 2°C to 6°C except as noted below.

General QA/QC issues associated with the analytical data enclosed in this laboratory report will be indicated with a reference to a comment or explanation on the Data Qualifier page. More complex and involved QA/QC issues will be discussed in detail below.

NWTPH Gx Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Volatiles (soil) EPA 8260B Analysis

Per EPA Method 5035A, samples were received by the laboratory in pre-weighed 40 mL VOA vials within 48 hours of sample collection. They were stored in a freezer at between -7°C and -20°C until extraction or analysis.

The results for Toluene in samples B-4_6, B-5_6 and B-5_7 may be the result of contaminated sample vials.

Interfering compounds in sample B-4_6 may be masking the presence of Acetone.

Internal Standard 1,4-Dichlorobenzene-d4 does not meet acceptance criteria and Surrogate Standard 4-Bromofluorobenzene is outside of the control limits for sample B-5_6 due to sample matrix effects. The sample was re-analyzed at a dilution with normal standard recoveries. All results, including Practical Quantitation Limits, from Bromobenzene onward should be considered estimates.

The values reported for Acetone in samples B-5_6 and B-5_7 exceed the quantitation range and are therefore estimates. The samples were re-analyzed at the lowest possible dilution allowed by Method 5035 with non-detect results for Acetone.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Project: 8915 162890

Case Narrative Page 2 of 2

Volatiles (water) EPA 8260B Analysis

Sample TRIP BLANK 2 was not analyzed for Vinyl Acetate.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Organochlorine Pesticides (soil) by EPA 8081A Analysis

Due to negative effects of the matrix on the instrument, values for the analytes 4,4'-DDT and Methoxychlor in the closing continuing calibration verification standard (CCV) were low. Therefore, values for these compounds can be greater than reported. Since the degradation of the CCV standards was reproducible after re-injecting the sample extracts, the CCV degradation problem was attributed to the matrix of these samples.

Any other QA/QC issues associated with this extraction and analysis will be indicated with a footnote reference and discussed in detail on the Data Qualifier page.

Project: 8915 162890

NWTPH-Gx

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Client ID: **B-4_6 B-5_6** Lab ID: 03-076-02 03-076-05

| | Result | Flags | PQL | Result | Flags | PQL |
|--------------------------------------|--------|-------|-----|--------|-------|-----|
| TPH-Gas | ND | | 4.6 | ND | | 6.0 |
| Surrogate Recovery: Fluorobenzene | 89% | | | 94% | | |

Project: 8915 162890

NWTPH-Gx

Date Extracted: 3-12-08
Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Client ID: **B-5_7**Lab ID: 03-076-06

Result Flags PQL

TPH-Gas ND 5.1

Surrogate Recovery:

Fluorobenzene 92%

Project: 8915 162890

NWTPH-Gx METHOD BLANK QUALITY CONTROL

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0312S1

Result Flags PQL

TPH-Gas ND 5.0

Surrogate Recovery:

Fluorobenzene 99%

Project: 8915 162890

NWTPH-Gx DUPLICATE QUALITY CONTROL

Date Extracted: 3-12-08
Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-01 03-062-01

Original Duplicate RPD Flags

TPH-Gas ND ND NA

Surrogate Recovery:

Fluorobenzene 94% 94%

Project: 8915 162890

NWTPH-Gx

Date Extracted: 3-13-08 Date Analyzed: 3-13-08

Matrix: Water Units: ug/L (ppb)

Client ID: EB-1

Lab ID: 03-076-13

Result Flags PQL

TPH-Gas ND 100

Surrogate Recovery:

Fluorobenzene 106%

Project: 8915 162890

NWTPH-Gx METHOD BLANK QUALITY CONTROL

Date Extracted: 3-13-08 Date Analyzed: 3-13-08

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0313W1

Result Flags PQL

TPH-Gas ND 100

Surrogate Recovery:

Fluorobenzene 107%

Project: 8915 162890

NWTPH-Gx DUPLICATE QUALITY CONTROL

Date Extracted: 3-13-08 Date Analyzed: 3-13-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-076-13 03-076-13

Original Duplicate RPD Flags

TPH-Gas ND ND NA

Surrogate Recovery:

Fluorobenzene 106% 104%

Project: 8915 162890

NWTPH-Dx

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

| Client ID: Lab ID: | B-4_6 03-076-02 | B-5_6 03-076-05 | B-5_7 03-076-06 |
|---------------------------------|---------------------------|---------------------------|---------------------------|
| Diesel Range: PQL: | ND 28 | ND 140 | ND 140 |
| Identification: | | | |
| Lube Oil Range: | 110 56 | 2700 280 | 2400 270 |
| Identification: | Lube Oil | Lube Oil | Lube Oil |
| Surrogate Recovery o-Terphenyl: | 71% | 79% | 79% |
| Flags: | Υ | Υ | Υ |

Project: 8915 162890

NWTPH-Dx METHOD BLANK QUALITY CONTROL

| Date Extracted: | 3-12-08 |
|-----------------|---------|
| Date Analyzed: | 3-12-08 |

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0312S1

Diesel Range: ND PQL: 25

Identification: ---

Lube Oil Range: ND PQL: 50

Identification: ---

Surrogate Recovery

o-Terphenyl: 88%

Flags: Y

Project: 8915 162890

NWTPH-Dx DUPLICATE QUALITY CONTROL

03-062-01 DUP

ND

25

| Date Extracted: | 3-12-08 |
|-----------------|-----------------|
| Date Analyzed: | 3-12-08 |
| Matrix: | Soil |
| Units: | mg/kg (ppm) |
| Lab ID: | 03-062-01 |
| Diesel Range: | ND 25 |

RPD: N/A

Surrogate Recovery

o-Terphenyl: 84% 75%

Flags: Y Y

Project: 8915 162890

NWTPH-Dx

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Water Units: mg/L (ppm)

Client ID: EB-1Lab ID: 03-076-13

Diesel Range: ND PQL: 0.30

Identification: ---

Lube Oil Range: **ND**PQL: 0.47

Identification: ---

Surrogate Recovery

o-Terphenyl: 82%

Flags: Y

Project: 8915 162890

NWTPH-Dx METHOD BLANK QUALITY CONTROL

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Water Units: mg/L (ppm)

Lab ID: MB0314W1

Diesel Range: **ND**PQL: 0.25

Identification: ---

Lube Oil Range: **ND**PQL: 0.40

Identification: ---

Surrogate Recovery

o-Terphenyl: 76%

Flags: Y

Project: 8915 162890

NWTPH-Dx DUPLICATE QUALITY CONTROL

| Date Extracted: | 3-14-08 |
|-----------------|---------|
| Date Analyzed: | 3-18-08 |

Matrix: Water Units: mg/L (ppm)

Lab ID: 03-092-01 03-092-01 DUP

Diesel Range: ND ND PQL: 0.26 0.25

RPD: N/A

Surrogate Recovery

o-Terphenyl: 82% 81%

Flags: Y Y

Project: 8915 162890

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-076-02 **Client ID: B-4_6**

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|---------|
| Dichlorodifluoromethane | ND | · · | 0.00095 |
| Chloromethane | ND | | 0.0047 |
| Vinyl Chloride | ND | | 0.00095 |
| Bromomethane | ND | | 0.00095 |
| Chloroethane | ND | | 0.0047 |
| Trichlorofluoromethane | ND | | 0.00095 |
| 1,1-Dichloroethene | ND | | 0.00095 |
| Acetone | ND | | 0.0047 |
| lodomethane | ND | | 0.0047 |
| Carbon Disulfide | ND | | 0.00095 |
| Methylene Chloride | ND | | 0.0047 |
| (trans) 1,2-Dichloroethene | ND | | 0.00095 |
| Methyl t-Butyl Ether | ND | | 0.00095 |
| 1,1-Dichloroethane | ND | | 0.00095 |
| Vinyl Acetate | ND | | 0.0047 |
| 2,2-Dichloropropane | ND | | 0.00095 |
| (cis) 1,2-Dichloroethene | ND | | 0.00095 |
| 2-Butanone | ND | | 0.0047 |
| Bromochloromethane | ND | | 0.00095 |
| Chloroform | ND | | 0.00095 |
| 1,1,1-Trichloroethane | ND | | 0.00095 |
| Carbon Tetrachloride | ND | | 0.00095 |
| 1,1-Dichloropropene | ND | | 0.00095 |
| Benzene | ND | | 0.00095 |
| 1,2-Dichloroethane | ND | | 0.00095 |
| Trichloroethene | ND | | 0.00095 |
| 1,2-Dichloropropane | ND | | 0.00095 |
| Dibromomethane | ND | | 0.00095 |
| Bromodichloromethane | ND | | 0.00095 |
| 2-Chloroethyl Vinyl Ether | ND | | 0.0047 |
| (cis) 1,3-Dichloropropene | ND | | 0.00095 |
| Methyl Isobutyl Ketone | ND | | 0.0047 |
| Toluene | 0.0029 | Z | 0.00095 |
| (trans) 1,3-Dichloropropene | ND | | 0.00095 |

Project: 8915 162890

VOLATILES by EPA 8260B Page 2 of 2

Lab ID: 03-076-02 B-4_6 Client ID:

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|---------|
| 1,1,2-Trichloroethane | ND | | 0.00095 |
| Tetrachloroethene | ND | | 0.00095 |
| 1,3-Dichloropropane | ND | | 0.00095 |
| 2-Hexanone | ND | | 0.0047 |
| Dibromochloromethane | ND | | 0.00095 |
| 1,2-Dibromoethane | ND | | 0.00095 |
| Chlorobenzene | ND | | 0.00095 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.00095 |
| Ethylbenzene | ND | | 0.00095 |
| m,p-Xylene | ND | | 0.0019 |
| o-Xylene | ND | | 0.00095 |
| Styrene | ND | | 0.00095 |
| Bromoform | ND | | 0.00095 |
| Isopropylbenzene | ND | | 0.00095 |
| Bromobenzene | ND | | 0.00095 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.00095 |
| 1,2,3-Trichloropropane | ND | | 0.00095 |
| n-Propylbenzene | ND | | 0.00095 |
| 2-Chlorotoluene | ND | | 0.00095 |
| 4-Chlorotoluene | ND | | 0.00095 |
| 1,3,5-Trimethylbenzene | ND | | 0.00095 |
| tert-Butylbenzene | ND | | 0.00095 |
| 1,2,4-Trimethylbenzene | ND | | 0.00095 |
| sec-Butylbenzene | ND | | 0.00095 |
| 1,3-Dichlorobenzene | ND | | 0.00095 |
| p-Isopropyltoluene | ND | | 0.00095 |
| 1,4-Dichlorobenzene | ND | | 0.00095 |
| 1,2-Dichlorobenzene | ND | | 0.00095 |
| n-Butylbenzene | ND | | 0.00095 |
| 1,2-Dibromo-3-chloropropane | ND | | 0.0047 |
| 1,2,4-Trichlorobenzene | ND | | 0.00095 |
| Hexachlorobutadiene | ND | | 0.0047 |
| Naphthalene | ND | | 0.00095 |
| 1,2,3-Trichlorobenzene | ND | | 0.00095 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 77 | 70-118 |
| Toluene-d8 | 85 | 70-121 |
| 4-Bromofluorobenzene | 86 | 70-130 |

Project: 8915 162890

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-076-05 **Client ID: B-5_6**

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|--------|
| Dichlorodifluoromethane | ND | · · | 0.0011 |
| Chloromethane | ND | | 0.0055 |
| Vinyl Chloride | ND | | 0.0011 |
| Bromomethane | ND | | 0.0011 |
| Chloroethane | ND | | 0.0055 |
| Trichlorofluoromethane | ND | | 0.0011 |
| 1,1-Dichloroethene | ND | | 0.0011 |
| Acetone | 0.73 | E | 0.0055 |
| lodomethane | ND | | 0.0055 |
| Carbon Disulfide | ND | | 0.0011 |
| Methylene Chloride | ND | | 0.0055 |
| (trans) 1,2-Dichloroethene | ND | | 0.0011 |
| Methyl t-Butyl Ether | ND | | 0.0011 |
| 1,1-Dichloroethane | ND | | 0.0011 |
| Vinyl Acetate | ND | | 0.0055 |
| 2,2-Dichloropropane | ND | | 0.0011 |
| (cis) 1,2-Dichloroethene | ND | | 0.0011 |
| 2-Butanone | ND | | 0.0055 |
| Bromochloromethane | ND | | 0.0011 |
| Chloroform | ND | | 0.0011 |
| 1,1,1-Trichloroethane | ND | | 0.0011 |
| Carbon Tetrachloride | ND | | 0.0011 |
| 1,1-Dichloropropene | ND | | 0.0011 |
| Benzene | ND | | 0.0011 |
| 1,2-Dichloroethane | ND | | 0.0011 |
| Trichloroethene | ND | | 0.0011 |
| 1,2-Dichloropropane | ND | | 0.0011 |
| Dibromomethane | ND | | 0.0011 |
| Bromodichloromethane | ND | | 0.0011 |
| 2-Chloroethyl Vinyl Ether | ND | | 0.0055 |
| (cis) 1,3-Dichloropropene | ND | | 0.0011 |
| Methyl Isobutyl Ketone | ND | | 0.0055 |
| Toluene | 0.0016 | Z | 0.0011 |
| (trans) 1,3-Dichloropropene | ND | | 0.0011 |

Project: 8915 162890

VOLATILES by EPA 8260B Page 2 of 2

Lab ID: 03-076-05 Client ID: B-5_6

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|--------|
| 1,1,2-Trichloroethane | ND | | 0.0011 |
| Tetrachloroethene | 0.0053 | | 0.0011 |
| 1,3-Dichloropropane | ND | | 0.0011 |
| 2-Hexanone | ND | | 0.0055 |
| Dibromochloromethane | ND | | 0.0011 |
| 1,2-Dibromoethane | ND | | 0.0011 |
| Chlorobenzene | ND | | 0.0011 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.0011 |
| Ethylbenzene | ND | | 0.0011 |
| m,p-Xylene | ND | | 0.0022 |
| o-Xylene | ND | | 0.0011 |
| Styrene | ND | | 0.0011 |
| Bromoform | ND | | 0.0011 |
| Isopropylbenzene | ND | | 0.0011 |
| Bromobenzene | ND | | 0.0011 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.0011 |
| 1,2,3-Trichloropropane | ND | | 0.0011 |
| n-Propylbenzene | ND | | 0.0011 |
| 2-Chlorotoluene | ND | | 0.0011 |
| 4-Chlorotoluene | ND | | 0.0011 |
| 1,3,5-Trimethylbenzene | ND | | 0.0011 |
| tert-Butylbenzene | ND | | 0.0011 |
| 1,2,4-Trimethylbenzene | ND | | 0.0011 |
| sec-Butylbenzene | ND | | 0.0011 |
| 1,3-Dichlorobenzene | ND | | 0.0011 |
| p-Isopropyltoluene | ND | | 0.0011 |
| 1,4-Dichlorobenzene | ND | | 0.0011 |
| 1,2-Dichlorobenzene | ND | | 0.0011 |
| n-Butylbenzene | ND | | 0.0011 |
| 1,2-Dibromo-3-chloropropane | ND | | 0.0055 |
| 1,2,4-Trichlorobenzene | ND | | 0.0011 |
| Hexachlorobutadiene | ND | | 0.0055 |
| Naphthalene | ND | | 0.0011 |
| 1,2,3-Trichlorobenzene | ND | | 0.0011 |

| | Percent | | Control |
|----------------------|----------|---|---------|
| Surrogate | Recovery | | Limits |
| Dibromofluoromethane | 81 | | 70-118 |
| Toluene-d8 | 81 | | 70-121 |
| 4-Bromofluorobenzene | 62 | Q | 70-130 |

Project: 8915 162890

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-076-06

Client ID: B-5_7

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|--------|
| Dichlorodifluoromethane | ND | | 0.0010 |
| Chloromethane | ND | | 0.0050 |
| Vinyl Chloride | ND | | 0.0010 |
| Bromomethane | ND | | 0.0010 |
| Chloroethane | ND | | 0.0050 |
| Trichlorofluoromethane | ND | | 0.0010 |
| 1,1-Dichloroethene | ND | | 0.0010 |
| Acetone | 0.38 | Е | 0.0050 |
| lodomethane | ND | | 0.0050 |
| Carbon Disulfide | ND | | 0.0010 |
| Methylene Chloride | ND | | 0.0050 |
| (trans) 1,2-Dichloroethene | ND | | 0.0010 |
| Methyl t-Butyl Ether | ND | | 0.0010 |
| 1,1-Dichloroethane | ND | | 0.0010 |
| Vinyl Acetate | ND | | 0.0050 |
| 2,2-Dichloropropane | ND | | 0.0010 |
| (cis) 1,2-Dichloroethene | ND | | 0.0010 |
| 2-Butanone | ND | | 0.0050 |
| Bromochloromethane | ND | | 0.0010 |
| Chloroform | ND | | 0.0010 |
| 1,1,1-Trichloroethane | ND | | 0.0010 |
| Carbon Tetrachloride | ND | | 0.0010 |
| 1,1-Dichloropropene | ND | | 0.0010 |
| Benzene | ND | | 0.0010 |
| 1,2-Dichloroethane | ND | | 0.0010 |
| Trichloroethene | ND | | 0.0010 |
| 1,2-Dichloropropane | ND | | 0.0010 |
| Dibromomethane | ND | | 0.0010 |
| Bromodichloromethane | ND | | 0.0010 |
| 2-Chloroethyl Vinyl Ether | ND | | 0.0050 |
| (cis) 1,3-Dichloropropene | ND | | 0.0010 |
| Methyl Isobutyl Ketone | ND | | 0.0050 |
| Toluene | 0.0026 | Z | 0.0010 |
| (trans) 1,3-Dichloropropene | ND | | 0.0010 |

Project: 8915 162890

VOLATILES by EPA 8260B Page 2 of 2

Lab ID: 03-076-06 Client ID: B-5_7

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|--------|
| 1,1,2-Trichloroethane | ND | J | 0.0010 |
| Tetrachloroethene | 0.0013 | | 0.0010 |
| 1,3-Dichloropropane | ND | | 0.0010 |
| 2-Hexanone | ND | | 0.0050 |
| Dibromochloromethane | ND | | 0.0010 |
| 1,2-Dibromoethane | ND | | 0.0010 |
| Chlorobenzene | ND | | 0.0010 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.0010 |
| Ethylbenzene | ND | | 0.0010 |
| m,p-Xylene | ND | | 0.0020 |
| o-Xylene | ND | | 0.0010 |
| Styrene | ND | | 0.0010 |
| Bromoform | ND | | 0.0010 |
| Isopropylbenzene | ND | | 0.0010 |
| Bromobenzene | ND | | 0.0010 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.0010 |
| 1,2,3-Trichloropropane | ND | | 0.0010 |
| n-Propylbenzene | ND | | 0.0010 |
| 2-Chlorotoluene | ND | | 0.0010 |
| 4-Chlorotoluene | ND | | 0.0010 |
| 1,3,5-Trimethylbenzene | ND | | 0.0010 |
| tert-Butylbenzene | ND | | 0.0010 |
| 1,2,4-Trimethylbenzene | ND | | 0.0010 |
| sec-Butylbenzene | ND | | 0.0010 |
| 1,3-Dichlorobenzene | ND | | 0.0010 |
| p-Isopropyltoluene | ND | | 0.0010 |
| 1,4-Dichlorobenzene | ND | | 0.0010 |
| 1,2-Dichlorobenzene | ND | | 0.0010 |
| n-Butylbenzene | ND | | 0.0010 |
| 1,2-Dibromo-3-chloropropane | ND | | 0.0050 |
| 1,2,4-Trichlorobenzene | ND | | 0.0010 |
| Hexachlorobutadiene | ND | | 0.0050 |
| Naphthalene | ND | | 0.0010 |
| 1,2,3-Trichlorobenzene | ND | | 0.0010 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 81 | 70-118 |
| Toluene-d8 | 88 | 70-121 |
| 4-Bromofluorobenzene | 76 | 70-130 |

Project: 8915 162890

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 1 of 2

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0314S1

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|--------|
| Dichlorodifluoromethane | ND | _ | 0.0010 |
| Chloromethane | ND | | 0.0050 |
| Vinyl Chloride | ND | | 0.0010 |
| Bromomethane | ND | | 0.0010 |
| Chloroethane | ND | | 0.0050 |
| Trichlorofluoromethane | ND | | 0.0010 |
| 1,1-Dichloroethene | ND | | 0.0010 |
| Acetone | ND | | 0.0050 |
| lodomethane | ND | | 0.0050 |
| Carbon Disulfide | ND | | 0.0010 |
| Methylene Chloride | ND | | 0.0050 |
| (trans) 1,2-Dichloroethene | ND | | 0.0010 |
| Methyl t-Butyl Ether | ND | | 0.0010 |
| 1,1-Dichloroethane | ND | | 0.0010 |
| Vinyl Acetate | ND | | 0.0050 |
| 2,2-Dichloropropane | ND | | 0.0010 |
| (cis) 1,2-Dichloroethene | ND | | 0.0010 |
| 2-Butanone | ND | | 0.0050 |
| Bromochloromethane | ND | | 0.0010 |
| Chloroform | ND | | 0.0010 |
| 1,1,1-Trichloroethane | ND | | 0.0010 |
| Carbon Tetrachloride | ND | | 0.0010 |
| 1,1-Dichloropropene | ND | | 0.0010 |
| Benzene | ND | | 0.0010 |
| 1,2-Dichloroethane | ND | | 0.0010 |
| Trichloroethene | ND | | 0.0010 |
| 1,2-Dichloropropane | ND | | 0.0010 |
| Dibromomethane | ND | | 0.0010 |
| Bromodichloromethane | ND | | 0.0010 |
| 2-Chloroethyl Vinyl Ether | ND | | 0.0050 |
| (cis) 1,3-Dichloropropene | ND | | 0.0010 |
| Methyl Isobutyl Ketone | ND | | 0.0050 |
| Toluene | ND | | 0.0010 |
| (trans) 1,3-Dichloropropene | ND | | 0.0010 |

Project: 8915 162890

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 2 of 2

Lab ID: MB0314S1

| Compound 1,1,2-Trichloroethane Tetrachloroethene 1,3-Dichloropropane 2-Hexanone Dibromochloromethane 1,2-Dibromoethane | ND N | Flags | PQL 0.0010 0.0010 0.0050 0.0010 0.0010 0.0010 0.0010 0.0020 0.0010 |
|---|--|-------|---|
| Tetrachloroethene 1,3-Dichloropropane 2-Hexanone Dibromochloromethane 1,2-Dibromoethane | ND ND ND ND ND ND ND ND | | 0.0010 0.0050 0.0010 0.0010 0.0010 0.0010 0.0020 0.0010 |
| 2-Hexanone Dibromochloromethane 1,2-Dibromoethane | ND ND ND ND ND ND ND ND | | 0.0050 0.0010 0.0010 0.0010 0.0010 0.0020 0.0010 |
| Dibromochloromethane 1,2-Dibromoethane | ND ND ND ND ND ND ND | | 0.0010 0.0010 0.0010 0.0010 0.0010 0.0020 0.0010 |
| 1,2-Dibromoethane | ND ND ND ND ND ND | | 0.0010 0.0010 0.0010 0.0010 0.0020 0.0010 |
| • | ND ND ND ND ND ND | | 0.0010 0.0010 0.0010 0.0020 0.0010 |
| Chlorobonzono | ND ND ND ND ND | | 0.0010 0.0010 0.0020 0.0010 |
| Chlorobenzene | ND ND ND ND | | 0.0010 0.0020 0.0010 |
| 1,1,1,2-Tetrachloroethane | ND ND ND | | 0.0020 0.0010 |
| Ethylbenzene | ND ND | | 0.0010 |
| m,p-Xylene | ND | | |
| o-Xylene | | | 0 0010 |
| Styrene | ND | | 0.0010 |
| Bromoform | | | 0.0010 |
| Isopropylbenzene | ND | | 0.0010 |
| Bromobenzene | ND | | 0.0010 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.0010 |
| 1,2,3-Trichloropropane | ND | | 0.0010 |
| n-Propylbenzene | ND | | 0.0010 |
| 2-Chlorotoluene | ND | | 0.0010 |
| 4-Chlorotoluene | ND | | 0.0010 |
| 1,3,5-Trimethylbenzene | ND | | 0.0010 |
| tert-Butylbenzene | ND | | 0.0010 |
| 1,2,4-Trimethylbenzene | ND | | 0.0010 |
| sec-Butylbenzene | ND | | 0.0010 |
| 1,3-Dichlorobenzene | ND | | 0.0010 |
| p-Isopropyltoluene | ND | | 0.0010 |
| 1,4-Dichlorobenzene | ND | | 0.0010 |
| 1,2-Dichlorobenzene | ND | | 0.0010 |
| n-Butylbenzene | ND | | 0.0010 |
| 1,2-Dibromo-3-chloropropane | ND | | 0.0050 |
| 1,2,4-Trichlorobenzene | ND | | 0.0010 |
| Hexachlorobutadiene | ND | | 0.0050 |
| Naphthalene | ND | | 0.0010 |
| 1,2,3-Trichlorobenzene | ND | | 0.0010 |
| Surregate | Percent | | Control |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 70 | 70-118 |
| Toluene-d8 | 91 | 70-121 |
| 4-Bromofluorobenzene | 79 | 70-130 |

Project: 8915 162890

VOLATILES by EPA 8260B SB/SBD QUALITY CONTROL

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: SB0314S1

| | Spike | | Percent | | Percent | Recovery | |
|--------------------|--------|--------|----------|--------|----------|----------|-------|
| Compound | Amount | SB | Recovery | SBD | Recovery | Limits | Flags |
| | | | | | | | |
| 1,1-Dichloroethene | 0.0500 | 0.0373 | 75 | 0.0412 | 82 | 70-130 | |
| Benzene | 0.0500 | 0.0419 | 84 | 0.0422 | 84 | 70-127 | |
| Trichloroethene | 0.0500 | 0.0404 | 81 | 0.0449 | 90 | 73-117 | |
| Toluene | 0.0500 | 0.0410 | 82 | 0.0436 | 87 | 78-115 | |
| Chlorobenzene | 0.0500 | 0.0410 | 82 | 0.0418 | 84 | 80-117 | |

| Limit | Flags |
|-------|----------------|
| | |
| 10 | |
| 11 | |
| 13 | |
| 12 | |
| 10 | |
| | 11 13 12 |

Project: 8915 162890

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-19-08 Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| Dichlorodifluoromethane | ND | | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | ND | | 0.20 |
| Acetone | ND | | 5.0 |
| lodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | 2.5 | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | ND | | 0.20 |
| Vinyl Acetate | ND | | 1.0 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | 1.7 | | 0.20 |
| 1,1,1-Trichloroethane | ND | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | ND | | 0.20 |
| 1,2-Dichloroethane | ND | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8915 162890

VOLATILES by EPA 8260B Page 2 of 2

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | | 0.20 |
| Tetrachloroethene | ND | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | ND | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | ND | | 0.20 |
| m,p-Xylene | ND | | 0.40 |
| o-Xylene | ND | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 88 | 71-126 |
| Toluene-d8 | 84 | 76-116 |
| 4-Bromofluorobenzene | 76 | 70-123 |

Project: 8915 162890

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 1 of 2

Date Extracted: 3-19-08 Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0319W1

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| Dichlorodifluoromethane | ND | _ | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | ND | | 0.20 |
| Acetone | ND | | 5.0 |
| Iodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | ND | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | ND | | 0.20 |
| Vinyl Acetate | ND | | 1.0 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | ND | | 0.20 |
| 1,1,1-Trichloroethane | ND | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | ND | | 0.20 |
| 1,2-Dichloroethane | ND | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8915 162890

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 2 of 2

Lab ID: MB0319W1

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | | 0.20 |
| Tetrachloroethene | ND | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | ND | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | ND | | 0.20 |
| m,p-Xylene | ND | | 0.40 |
| o-Xylene | ND | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 88 | 71-126 |
| Toluene-d8 | 83 | 76-116 |
| 4-Bromofluorobenzene | 76 | 70-123 |

Project: 8915 162890

VOLATILES by EPA 8260B SB/SBD QUALITY CONTROL

Date Extracted: 3-19-08 Date Analyzed: 3-19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: SB0319W1

| Compound | Spike Amount | SB | Percent Recovery | SBD | Percent Recovery | Recovery Limits | Flags |
|--------------------|-----------------|------|---------------------|------|---------------------|--------------------|-------|
| 1,1-Dichloroethene | 10.0 | 8.91 | 89 | 8.16 | 82 | 70-130 | |
| Benzene | 10.0 | 9.55 | 96 | 9.55 | 96 | 70-130 | |
| Trichloroethene | 10.0 | 8.45 | 85 | 8.29 | 83 | 70-116 | |
| Toluene | 10.0 | 9.80 | 98 | 9.62 | 96 | 76-119 | |
| Chlorobenzene | 10.0 | 9.12 | 91 | 9.16 | 92 | 77-112 | |

| | | RPD | |
|--------------------|-----|-------|-------|
| | RPD | Limit | Flags |
| 1,1-Dichloroethene | 9 | 20 | |
| Benzene | 0 | 16 | |
| Trichloroethene | 2 | 16 | |
| Toluene | 2 | 15 | |
| Chlorobenzene | 0 | 15 | |

Project: 8915 162890

VOLATILES by EPA 8260B

Page 1 of 2

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-076-11

Client ID: TRIP BLANK 2

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| Dichlorodifluoromethane | ND | • | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | ND | | 0.20 |
| Acetone | ND | | 5.0 |
| lodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | ND | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | ND | | 0.20 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | ND | | 0.20 |
| 1,1,1-Trichloroethane | ND | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | ND | | 0.20 |
| 1,2-Dichloroethane | ND | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8915 162890

VOLATILES by EPA 8260B Page 2 of 2

Lab ID: 03-076-11 **Client ID: TRIP BLANK 2**

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | | 0.20 |
| Tetrachloroethene | ND | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | ND | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | ND | | 0.20 |
| m,p-Xylene | ND | | 0.40 |
| o-Xylene | ND | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 92 | 71-126 |
| Toluene-d8 | 86 | 76-116 |
| 4-Bromofluorobenzene | 83 | 70-123 |

Project: 8915 162890

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 1 of 2

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0314W1

| Compound | Results | Flags | PQL |
|-----------------------------|---------|--------|------|
| Dichlorodifluoromethane | ND | i lago | 0.20 |
| Chloromethane | ND | | 1.0 |
| Vinyl Chloride | ND | | 0.20 |
| Bromomethane | ND | | 0.20 |
| Chloroethane | ND | | 1.0 |
| Trichlorofluoromethane | ND | | 0.20 |
| 1,1-Dichloroethene | ND | | 0.20 |
| Acetone | ND | | 5.0 |
| lodomethane | ND | | 1.0 |
| Carbon Disulfide | ND | | 0.20 |
| Methylene Chloride | ND | | 1.0 |
| (trans) 1,2-Dichloroethene | ND | | 0.20 |
| Methyl t-Butyl Ether | ND | | 0.20 |
| 1,1-Dichloroethane | ND | | 0.20 |
| 2,2-Dichloropropane | ND | | 0.20 |
| (cis) 1,2-Dichloroethene | ND | | 0.20 |
| 2-Butanone | ND | | 5.0 |
| Bromochloromethane | ND | | 0.20 |
| Chloroform | ND | | 0.20 |
| 1,1,1-Trichloroethane | ND | | 0.20 |
| Carbon Tetrachloride | ND | | 0.20 |
| 1,1-Dichloropropene | ND | | 0.20 |
| Benzene | ND | | 0.20 |
| 1,2-Dichloroethane | ND | | 0.20 |
| Trichloroethene | ND | | 0.20 |
| 1,2-Dichloropropane | ND | | 0.20 |
| Dibromomethane | ND | | 0.20 |
| Bromodichloromethane | ND | | 0.20 |
| 2-Chloroethyl Vinyl Ether | ND | | 1.0 |
| (cis) 1,3-Dichloropropene | ND | | 0.20 |
| Methyl Isobutyl Ketone | ND | | 2.0 |
| Toluene | ND | | 0.20 |
| (trans) 1,3-Dichloropropene | ND | | 0.20 |

Project: 8915 162890

VOLATILES by EPA 8260B METHOD BLANK QUALITY CONTROL

Page 2 of 2

Lab ID: MB0314W1

| Compound | Results | Flags | PQL |
|-----------------------------|---------|-------|------|
| 1,1,2-Trichloroethane | ND | | 0.20 |
| Tetrachloroethene | ND | | 0.20 |
| 1,3-Dichloropropane | ND | | 0.20 |
| 2-Hexanone | ND | | 2.0 |
| Dibromochloromethane | ND | | 0.20 |
| 1,2-Dibromoethane | ND | | 0.20 |
| Chlorobenzene | ND | | 0.20 |
| 1,1,1,2-Tetrachloroethane | ND | | 0.20 |
| Ethylbenzene | ND | | 0.20 |
| m,p-Xylene | ND | | 0.40 |
| o-Xylene | ND | | 0.20 |
| Styrene | ND | | 0.20 |
| Bromoform | ND | | 1.0 |
| Isopropylbenzene | ND | | 0.20 |
| Bromobenzene | ND | | 0.20 |
| 1,1,2,2-Tetrachloroethane | ND | | 0.20 |
| 1,2,3-Trichloropropane | ND | | 0.20 |
| n-Propylbenzene | ND | | 0.20 |
| 2-Chlorotoluene | ND | | 0.20 |
| 4-Chlorotoluene | ND | | 0.20 |
| 1,3,5-Trimethylbenzene | ND | | 0.20 |
| tert-Butylbenzene | ND | | 0.20 |
| 1,2,4-Trimethylbenzene | ND | | 0.20 |
| sec-Butylbenzene | ND | | 0.20 |
| 1,3-Dichlorobenzene | ND | | 0.20 |
| p-Isopropyltoluene | ND | | 0.20 |
| 1,4-Dichlorobenzene | ND | | 0.20 |
| 1,2-Dichlorobenzene | ND | | 0.20 |
| n-Butylbenzene | ND | | 0.20 |
| 1,2-Dibromo-3-chloropropane | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 0.20 |
| Hexachlorobutadiene | ND | | 0.20 |
| Naphthalene | ND | | 1.0 |
| 1,2,3-Trichlorobenzene | ND | | 0.20 |

| | Percent | Control |
|----------------------|----------|---------|
| Surrogate | Recovery | Limits |
| Dibromofluoromethane | 89 | 71-126 |
| Toluene-d8 | 85 | 76-116 |
| 4-Bromofluorobenzene | 83 | 70-123 |

Project: 8915 162890

VOLATILES by EPA 8260B SB/SBD QUALITY CONTROL

Date Extracted: 3-14-08 Date Analyzed: 3-14-08

Matrix: Water Units: ug/L (ppb)

Lab ID: SB0314W1

| | Spike | | Percent | | Percent | Recovery | |
|--------------------|--------|------|----------|------|----------|----------|-------|
| Compound | Amount | SB | Recovery | SBD | Recovery | Limits | Flags |
| 1,1-Dichloroethene | 10.0 | 11.4 | 114 | 10.6 | 106 | 70-130 | |
| Benzene | 10.0 | 10.8 | 108 | 11.1 | 111 | 70-130 | |
| Trichloroethene | 10.0 | 9.33 | 93 | 9.55 | 96 | 70-116 | |
| Toluene | 10.0 | 10.4 | 104 | 10.6 | 106 | 76-119 | |
| Chlorobenzene | 10.0 | 9.42 | 94 | 9.73 | 97 | 77-112 | |
| | | | | | | | |

| | RPD | | |
|--------------------|-----|-------|-------|
| | RPD | Limit | Flags |
| 1,1-Dichloroethene | 8 | 20 | |
| Benzene | 2 | 16 | |
| Trichloroethene | 2 | 16 | |
| Toluene | 2 | 15 | |
| Chlorobenzene | 3 | 15 | |
| | | | |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 1 of 3

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|--------|
| N-Nitrosodimethylamine | ND | • | 0.037 |
| Pyridine | ND | | 0.037 |
| Phenol | ND | | 0.037 |
| Aniline | ND | | 0.037 |
| bis(2-Chloroethyl)ether | ND | | 0.037 |
| 2-Chlorophenol | ND | | 0.037 |
| 1,3-Dichlorobenzene | ND | | 0.037 |
| 1,4-Dichlorobenzene | ND | | 0.037 |
| Benzyl alcohol | ND | | 0.037 |
| 1,2-Dichlorobenzene | ND | | 0.037 |
| 2-Methylphenol (o-Cresol) | ND | | 0.037 |
| bis(2-Chloroisopropyl)ether | ND | | 0.037 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 0.037 |
| N-Nitroso-di-n-propylamine | ND | | 0.037 |
| Hexachloroethane | ND | | 0.037 |
| Nitrobenzene | ND | | 0.037 |
| Isophorone | ND | | 0.037 |
| 2-Nitrophenol | ND | | 0.037 |
| 2,4-Dimethylphenol | ND | | 0.037 |
| bis(2-Chloroethoxy)methane | ND | | 0.037 |
| 2,4-Dichlorophenol | ND | | 0.037 |
| 1,2,4-Trichlorobenzene | ND | | 0.037 |
| Naphthalene | ND | | 0.0073 |
| 4-Chloroaniline | ND | | 0.037 |
| Hexachlorobutadiene | ND | | 0.037 |
| 4-Chloro-3-methylphenol | ND | | 0.037 |
| 2-Methylnaphthalene | ND | | 0.0073 |
| 1-Methylnaphthalene | ND | | 0.0073 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Hexachlorocyclopentadiene | ND | J | 0.037 |
| 2,4,6-Trichlorophenol | ND | | 0.037 |
| 2,3-Dichloroaniline | ND | | 0.037 |
| 2,4,5-Trichlorophenol | ND | | 0.037 |
| 2-Chloronaphthalene | ND | | 0.037 |
| 2-Nitroaniline | ND | | 0.037 |
| 1,4-Dinitrobenzene | ND | | 0.037 |
| Dimethylphthalate | ND | | 0.037 |
| 1,3-Dinitrobenzene | ND | | 0.037 |
| 2,6-Dinitrotoluene | ND | | 0.037 |
| 1,2-Dinitrobenzene | ND | | 0.037 |
| Acenaphthylene | ND | | 0.0073 |
| 3-Nitroaniline | ND | | 0.037 |
| 2,4-Dinitrophenol | ND | | 0.18 |
| Acenaphthene | ND | | 0.0073 |
| 4-Nitrophenol | ND | | 0.037 |
| 2,4-Dinitrotoluene | ND | | 0.037 |
| Dibenzofuran | ND | | 0.037 |
| 2,3,4,6-Tetrachlorophenol | ND | | 0.037 |
| 2,3,5,6-Tetrachlorophenol | ND | | 0.037 |
| Diethylphthalate | ND | | 0.037 |
| 4-Chlorophenyl-phenylether | ND | | 0.037 |
| 4-Nitroaniline | ND | | 0.037 |
| Fluorene | ND | | 0.0073 |
| 4,6-Dinitro-2-methylphenol | ND | | 0.18 |
| N-Nitrosodiphenylamine | ND | | 0.037 |
| 1,2-Diphenylhydrazine | ND | | 0.037 |
| 4-Bromophenyl-phenylether | ND | | 0.037 |
| Hexachlorobenzene | ND | | 0.037 |
| Pentachlorophenol | ND | | 0.18 |
| Phenanthrene | ND | | 0.0073 |
| Anthracene | ND | | 0.0073 |
| Carbazole | ND | | 0.037 |
| Di-n-butylphthalate | ND | | 0.037 |
| Fluoranthene | ND | | 0.0073 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

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| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Benzidine | ND | | 0.37 |
| Pyrene | ND | | 0.0073 |
| Butylbenzylphthalate | ND | | 0.037 |
| bis-2-Ethylhexyladipate | ND | | 0.037 |
| 3,3'-Dichlorobenzidine | ND | | 0.037 |
| Benzo[a]anthracene | ND | | 0.0073 |
| Chrysene | ND | | 0.0073 |
| bis(2-Ethylhexyl)phthalate | ND | | 0.037 |
| Di-n-octylphthalate | ND | | 0.037 |
| Benzo[b]fluoranthene | ND | | 0.0073 |
| Benzo[k]fluoranthene | ND | | 0.0073 |
| Benzo[a]pyrene | ND | | 0.0073 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.0073 |
| Dibenz[a,h]anthracene | ND | | 0.0073 |
| Benzo[g,h,i]perylene | ND | | 0.0073 |

| Surrogate : | Percent | Control |
|----------------------|----------|----------|
| _ | Recovery | Limits |
| 2-Fluorophenol | 81 | 39 - 90 |
| Phenol-d6 | 97 | 40 - 100 |
| Nitrobenzene-d5 | 84 | 30 - 100 |
| 2-Fluorobiphenyl | 84 | 41 - 100 |
| 2,4,6-Tribromophenol | 96 | 53 - 105 |
| Terphenyl-d14 | 105 | 49 - 115 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

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Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-076-07 **Client ID: B-5_8**

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|--------|
| N-Nitrosodimethylamine | ND | | 0.035 |
| Pyridine | ND | | 0.035 |
| Phenol | ND | | 0.035 |
| Aniline | ND | | 0.035 |
| bis(2-Chloroethyl)ether | ND | | 0.035 |
| 2-Chlorophenol | ND | | 0.035 |
| 1,3-Dichlorobenzene | ND | | 0.035 |
| 1,4-Dichlorobenzene | ND | | 0.035 |
| Benzyl alcohol | ND | | 0.035 |
| 1,2-Dichlorobenzene | ND | | 0.035 |
| 2-Methylphenol (o-Cresol) | ND | | 0.035 |
| bis(2-Chloroisopropyl)ether | ND | | 0.035 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 0.035 |
| N-Nitroso-di-n-propylamine | ND | | 0.035 |
| Hexachloroethane | ND | | 0.035 |
| Nitrobenzene | ND | | 0.035 |
| Isophorone | ND | | 0.035 |
| 2-Nitrophenol | ND | | 0.035 |
| 2,4-Dimethylphenol | ND | | 0.035 |
| bis(2-Chloroethoxy)methane | ND | | 0.035 |
| 2,4-Dichlorophenol | ND | | 0.035 |
| 1,2,4-Trichlorobenzene | ND | | 0.035 |
| Naphthalene | ND | | 0.0069 |
| 4-Chloroaniline | ND | | 0.035 |
| Hexachlorobutadiene | ND | | 0.035 |
| 4-Chloro-3-methylphenol | ND | | 0.035 |
| 2-Methylnaphthalene | ND | | 0.0069 |
| 1-Methylnaphthalene | ND | | 0.0069 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

Lab ID: 03-076-07 **Client ID: B-5_8**

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Hexachlorocyclopentadiene | ND | · · | 0.035 |
| 2,4,6-Trichlorophenol | ND | | 0.035 |
| 2,3-Dichloroaniline | ND | | 0.035 |
| 2,4,5-Trichlorophenol | ND | | 0.035 |
| 2-Chloronaphthalene | ND | | 0.035 |
| 2-Nitroaniline | ND | | 0.035 |
| 1,4-Dinitrobenzene | ND | | 0.035 |
| Dimethylphthalate | ND | | 0.035 |
| 1,3-Dinitrobenzene | ND | | 0.035 |
| 2,6-Dinitrotoluene | ND | | 0.035 |
| 1,2-Dinitrobenzene | ND | | 0.035 |
| Acenaphthylene | ND | | 0.0069 |
| 3-Nitroaniline | ND | | 0.035 |
| 2,4-Dinitrophenol | ND | | 0.17 |
| Acenaphthene | ND | | 0.0069 |
| 4-Nitrophenol | ND | | 0.035 |
| 2,4-Dinitrotoluene | ND | | 0.035 |
| Dibenzofuran | ND | | 0.035 |
| 2,3,4,6-Tetrachlorophenol | ND | | 0.035 |
| 2,3,5,6-Tetrachlorophenol | ND | | 0.035 |
| Diethylphthalate | ND | | 0.035 |
| 4-Chlorophenyl-phenylether | ND | | 0.035 |
| 4-Nitroaniline | ND | | 0.035 |
| Fluorene | ND | | 0.0069 |
| 4,6-Dinitro-2-methylphenol | ND | | 0.17 |
| N-Nitrosodiphenylamine | ND | | 0.035 |
| 1,2-Diphenylhydrazine | ND | | 0.035 |
| 4-Bromophenyl-phenylether | ND | | 0.035 |
| Hexachlorobenzene | ND | | 0.035 |
| Pentachlorophenol | ND | | 0.17 |
| Phenanthrene | ND | | 0.0069 |
| Anthracene | ND | | 0.0069 |
| Carbazole | ND | | 0.035 |
| Di-n-butylphthalate | ND | | 0.035 |
| Fluoranthene | ND | | 0.0069 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 3 of 3

Lab ID: 03-076-07 **Client ID: B-5_8**

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Benzidine | ND | | 0.35 |
| Pyrene | ND | | 0.0069 |
| Butylbenzylphthalate | ND | | 0.035 |
| bis-2-Ethylhexyladipate | ND | | 0.035 |
| 3,3'-Dichlorobenzidine | ND | | 0.035 |
| Benzo[a]anthracene | ND | | 0.0069 |
| Chrysene | ND | | 0.0069 |
| bis(2-Ethylhexyl)phthalate | ND | | 0.035 |
| Di-n-octylphthalate | ND | | 0.035 |
| Benzo[b]fluoranthene | ND | | 0.0069 |
| Benzo[k]fluoranthene | ND | | 0.0069 |
| Benzo[a]pyrene | ND | | 0.0069 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.0069 |
| Dibenz[a,h]anthracene | ND | | 0.0069 |
| Benzo[g,h,i]perylene | ND | | 0.0069 |

| Surrogate : | Percent | Control |
|----------------------|----------|----------|
| _ | Recovery | Limits |
| 2-Fluorophenol | 81 | 39 - 90 |
| Phenol-d6 | 96 | 40 - 100 |
| Nitrobenzene-d5 | 84 | 30 - 100 |
| 2-Fluorobiphenyl | 86 | 41 - 100 |
| 2,4,6-Tribromophenol | 93 | 53 - 105 |
| Terphenyl-d14 | 107 | 49 - 115 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 1 of 3

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-076-08 **Client ID:** 8-5_9

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|--------|
| N-Nitrosodimethylamine | ND | • | 0.035 |
| Pyridine | ND | | 0.035 |
| Phenol | ND | | 0.035 |
| Aniline | ND | | 0.035 |
| bis(2-Chloroethyl)ether | ND | | 0.035 |
| 2-Chlorophenol | ND | | 0.035 |
| 1,3-Dichlorobenzene | ND | | 0.035 |
| 1,4-Dichlorobenzene | ND | | 0.035 |
| Benzyl alcohol | ND | | 0.035 |
| 1,2-Dichlorobenzene | ND | | 0.035 |
| 2-Methylphenol (o-Cresol) | ND | | 0.035 |
| bis(2-Chloroisopropyl)ether | ND | | 0.035 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 0.035 |
| N-Nitroso-di-n-propylamine | ND | | 0.035 |
| Hexachloroethane | ND | | 0.035 |
| Nitrobenzene | ND | | 0.035 |
| Isophorone | ND | | 0.035 |
| 2-Nitrophenol | ND | | 0.035 |
| 2,4-Dimethylphenol | ND | | 0.035 |
| bis(2-Chloroethoxy)methane | ND | | 0.035 |
| 2,4-Dichlorophenol | ND | | 0.035 |
| 1,2,4-Trichlorobenzene | ND | | 0.035 |
| Naphthalene | ND | | 0.0069 |
| 4-Chloroaniline | ND | | 0.035 |
| Hexachlorobutadiene | ND | | 0.035 |
| 4-Chloro-3-methylphenol | ND | | 0.035 |
| 2-Methylnaphthalene | ND | | 0.0069 |
| 1-Methylnaphthalene | ND | | 0.0069 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

Lab ID: 03-076-08 **Client ID: B-5_9**

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Hexachlorocyclopentadiene | ND | J | 0.035 |
| 2,4,6-Trichlorophenol | ND | | 0.035 |
| 2,3-Dichloroaniline | ND | | 0.035 |
| 2,4,5-Trichlorophenol | ND | | 0.035 |
| 2-Chloronaphthalene | ND | | 0.035 |
| 2-Nitroaniline | ND | | 0.035 |
| 1,4-Dinitrobenzene | ND | | 0.035 |
| Dimethylphthalate | ND | | 0.035 |
| 1,3-Dinitrobenzene | ND | | 0.035 |
| 2,6-Dinitrotoluene | ND | | 0.035 |
| 1,2-Dinitrobenzene | ND | | 0.035 |
| Acenaphthylene | ND | | 0.0069 |
| 3-Nitroaniline | ND | | 0.035 |
| 2,4-Dinitrophenol | ND | | 0.17 |
| Acenaphthene | ND | | 0.0069 |
| 4-Nitrophenol | ND | | 0.035 |
| 2,4-Dinitrotoluene | ND | | 0.035 |
| Dibenzofuran | ND | | 0.035 |
| 2,3,4,6-Tetrachlorophenol | ND | | 0.035 |
| 2,3,5,6-Tetrachlorophenol | ND | | 0.035 |
| Diethylphthalate | ND | | 0.035 |
| 4-Chlorophenyl-phenylether | ND | | 0.035 |
| 4-Nitroaniline | ND | | 0.035 |
| Fluorene | ND | | 0.0069 |
| 4,6-Dinitro-2-methylphenol | ND | | 0.17 |
| N-Nitrosodiphenylamine | ND | | 0.035 |
| 1,2-Diphenylhydrazine | ND | | 0.035 |
| 4-Bromophenyl-phenylether | ND | | 0.035 |
| Hexachlorobenzene | ND | | 0.035 |
| Pentachlorophenol | ND | | 0.17 |
| Phenanthrene | ND | | 0.0069 |
| Anthracene | ND | | 0.0069 |
| Carbazole | ND | | 0.035 |
| Di-n-butylphthalate | 0.037 | | 0.035 |
| Fluoranthene | ND | | 0.0069 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 3 of 3

Lab ID: 03-076-08 **Client ID: B-5_9**

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Benzidine | ND | | 0.35 |
| Pyrene | ND | | 0.0069 |
| Butylbenzylphthalate | ND | | 0.035 |
| bis-2-Ethylhexyladipate | ND | | 0.035 |
| 3,3'-Dichlorobenzidine | ND | | 0.035 |
| Benzo[a]anthracene | ND | | 0.0069 |
| Chrysene | ND | | 0.0069 |
| bis(2-Ethylhexyl)phthalate | ND | | 0.035 |
| Di-n-octylphthalate | ND | | 0.035 |
| Benzo[b]fluoranthene | ND | | 0.0069 |
| Benzo[k]fluoranthene | ND | | 0.0069 |
| Benzo[a]pyrene | ND | | 0.0069 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.0069 |
| Dibenz[a,h]anthracene | ND | | 0.0069 |
| Benzo[g,h,i]perylene | ND | | 0.0069 |

| Surrogate: | Percent Recovery | Control Limits |
|----------------------|---------------------|-------------------|
| 2-Fluorophenol | 82 | 39 - 90 |
| Phenol-d6 | 96 | 40 - 100 |
| Nitrobenzene-d5 | 86 | 30 - 100 |
| 2-Fluorobiphenyl | 91 | 41 - 100 |
| 2,4,6-Tribromophenol | 98 | 53 - 105 |
| Terphenyl-d14 | 115 | 49 - 115 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 1 of 3

Date Extracted: 3-12-08 Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0312S1

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|--------|
| N-Nitrosodimethylamine | ND | | 0.033 |
| Pyridine | ND | | 0.033 |
| Phenol | ND | | 0.033 |
| Aniline | ND | | 0.033 |
| bis(2-Chloroethyl)ether | ND | | 0.033 |
| 2-Chlorophenol | ND | | 0.033 |
| 1,3-Dichlorobenzene | ND | | 0.033 |
| 1,4-Dichlorobenzene | ND | | 0.033 |
| Benzyl alcohol | ND | | 0.033 |
| 1,2-Dichlorobenzene | ND | | 0.033 |
| 2-Methylphenol (o-Cresol) | ND | | 0.033 |
| bis(2-Chloroisopropyl)ether | ND | | 0.033 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 0.033 |
| N-Nitroso-di-n-propylamine | ND | | 0.033 |
| Hexachloroethane | ND | | 0.033 |
| Nitrobenzene | ND | | 0.033 |
| Isophorone | ND | | 0.033 |
| 2-Nitrophenol | ND | | 0.033 |
| 2,4-Dimethylphenol | ND | | 0.033 |
| bis(2-Chloroethoxy)methane | ND | | 0.033 |
| 2,4-Dichlorophenol | ND | | 0.033 |
| 1,2,4-Trichlorobenzene | ND | | 0.033 |
| Naphthalene | ND | | 0.0067 |
| 4-Chloroaniline | ND | | 0.033 |
| Hexachlorobutadiene | ND | | 0.033 |
| 4-Chloro-3-methylphenol | ND | | 0.033 |
| 2-Methylnaphthalene | ND | | 0.0067 |
| 1-Methylnaphthalene | ND | | 0.0067 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 2 of 3

Lab ID: MB0312S1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Hexachlorocyclopentadiene | ND | _ | 0.033 |
| 2,4,6-Trichlorophenol | ND | | 0.033 |
| 2,3-Dichloroaniline | ND | | 0.033 |
| 2,4,5-Trichlorophenol | ND | | 0.033 |
| 2-Chloronaphthalene | ND | | 0.033 |
| 2-Nitroaniline | ND | | 0.033 |
| 1,4-Dinitrobenzene | ND | | 0.033 |
| Dimethylphthalate | ND | | 0.033 |
| 1,3-Dinitrobenzene | ND | | 0.033 |
| 2,6-Dinitrotoluene | ND | | 0.033 |
| 1,2-Dinitrobenzene | ND | | 0.033 |
| Acenaphthylene | ND | | 0.0067 |
| 3-Nitroaniline | ND | | 0.033 |
| 2,4-Dinitrophenol | ND | | 0.17 |
| Acenaphthene | ND | | 0.0067 |
| 4-Nitrophenol | ND | | 0.033 |
| 2,4-Dinitrotoluene | ND | | 0.033 |
| Dibenzofuran | ND | | 0.033 |
| 2,3,4,6-Tetrachlorophenol | ND | | 0.033 |
| 2,3,5,6-Tetrachlorophenol | ND | | 0.033 |
| Diethylphthalate | ND | | 0.033 |
| 4-Chlorophenyl-phenylether | ND | | 0.033 |
| 4-Nitroaniline | ND | | 0.033 |
| Fluorene | ND | | 0.0067 |
| 4,6-Dinitro-2-methylphenol | ND | | 0.17 |
| N-Nitrosodiphenylamine | ND | | 0.033 |
| 1,2-Diphenylhydrazine | ND | | 0.033 |
| 4-Bromophenyl-phenylether | ND | | 0.033 |
| Hexachlorobenzene | ND | | 0.033 |
| Pentachlorophenol | ND | | 0.17 |
| Phenanthrene | ND | | 0.0067 |
| Anthracene | ND | | 0.0067 |
| Carbazole | ND | | 0.033 |
| Di-n-butylphthalate | ND | | 0.033 |
| Fluoranthene | ND | | 0.0067 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 3 of 3

Lab ID: MB0312S1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|--------|
| Benzidine | ND | | 0.33 |
| Pyrene | ND | | 0.0067 |
| Butylbenzylphthalate | ND | | 0.033 |
| bis-2-Ethylhexyladipate | ND | | 0.033 |
| 3,3'-Dichlorobenzidine | ND | | 0.033 |
| Benzo[a]anthracene | ND | | 0.0067 |
| Chrysene | ND | | 0.0067 |
| bis(2-Ethylhexyl)phthalate | ND | | 0.033 |
| Di-n-octylphthalate | ND | | 0.033 |
| Benzo[b]fluoranthene | ND | | 0.0067 |
| Benzo[k]fluoranthene | ND | | 0.0067 |
| Benzo[a]pyrene | ND | | 0.0067 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.0067 |
| Dibenz[a,h]anthracene | ND | | 0.0067 |
| Benzo[g,h,i]perylene | ND | | 0.0067 |

| Surrogate : | Percent | Control |
|----------------------|----------|----------|
| | Recovery | Limits |
| 2-Fluorophenol | 76 | 39 - 90 |
| Phenol-d6 | 90 | 40 - 100 |
| Nitrobenzene-d5 | 84 | 30 - 100 |
| 2-Fluorobiphenyl | 87 | 41 - 100 |
| 2,4,6-Tribromophenol | 90 | 53 - 105 |
| Terphenyl-d14 | 103 | 49 - 115 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM SB/SBD QUALITY CONTROL

Date Extracted: 3-12-08
Date Analyzed: 3-12-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: SB0312S1

| | Spike | | Percent | | Percent | Recovery | • |
|----------------------------|--------|-------|----------|-------|----------|----------|-------|
| Compound: | Amount | SB | Recovery | SBD | Recovery | Limits | Flags |
| Phenol | 1.33 | 1.12 | 84 | 0.993 | 75 | 36-97 | |
| 2-Chlorophenol | 1.33 | 1.10 | 83 | 0.950 | 71 | 32-100 | |
| 1,4-Dichlorobenzene | 0.667 | 0.450 | 67 | 0.382 | 57 | 24-94 | |
| N-Nitroso-di-n-propylamine | 0.667 | 0.567 | 85 | 0.483 | 72 | 34-99 | |
| 1,2,4-Trichlorobenzene | 0.667 | 0.429 | 64 | 0.368 | 55 | 23-85 | |
| 4-Chloro-3-methylphenol | 1.33 | 1.18 | 89 | 1.18 | 88 | 46-108 | |
| Acenaphthene | 0.667 | 0.527 | 79 | 0.499 | 75 | 37-101 | |
| 2,4-Dinitrotoluene | 0.667 | 0.605 | 91 | 0.606 | 91 | 41-116 | |
| 4-Nitrophenol | 1.33 | 1.44 | 108 | 1.50 | 113 | 48-116 | |
| Pentachlorophenol | 1.33 | 1.21 | 91 | 1.25 | 94 | 28-130 | |
| Pyrene | 0.667 | 0.631 | 95 | 0.649 | 97 | 46-117 | |

| | | RPD | |
|----------------------------|-----|--------|-------|
| | RPD | Limits | Flags |
| Phenol | 12 | 29 | |
| 2-Chlorophenol | 14 | 28 | |
| 1,4-Dichlorobenzene | 16 | 27 | |
| N-Nitroso-di-n-propylamine | 16 | 30 | |
| 1,2,4-Trichlorobenzene | 15 | 28 | |
| 4-Chloro-3-methylphenol | 0 | 22 | |
| Acenaphthene | 6 | 25 | |
| 2,4-Dinitrotoluene | 0 | 30 | |
| 4-Nitrophenol | 4 | 30 | |
| Pentachlorophenol | 3 | 30 | |
| Pyrene | 3 | 20 | |
| | | | |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 1 of 3

Date Extracted: 3-13-08
Date Analyzed: 3-17&18-08

Matrix: Water Units: ug/L (ppb)

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|------|
| N-Nitrosodimethylamine | ND | • | 1.1 |
| Pyridine | ND | | 1.1 |
| Phenol | ND | | 1.1 |
| Aniline | ND | | 1.1 |
| bis(2-Chloroethyl)ether | ND | | 1.1 |
| 2-Chlorophenol | ND | | 1.1 |
| 1,3-Dichlorobenzene | ND | | 1.1 |
| 1,4-Dichlorobenzene | ND | | 1.1 |
| Benzyl alcohol | ND | | 1.1 |
| 1,2-Dichlorobenzene | ND | | 1.1 |
| 2-Methylphenol (o-Cresol) | ND | | 1.1 |
| bis(2-Chloroisopropyl)ether | ND | | 1.1 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 1.1 |
| N-Nitroso-di-n-propylamine | ND | | 1.1 |
| Hexachloroethane | ND | | 1.1 |
| Nitrobenzene | ND | | 1.1 |
| Isophorone | ND | | 1.1 |
| 2-Nitrophenol | ND | | 1.1 |
| 2,4-Dimethylphenol | ND | | 1.1 |
| bis(2-Chloroethoxy)methane | ND | | 1.1 |
| 2,4-Dichlorophenol | ND | | 1.1 |
| 1,2,4-Trichlorobenzene | ND | | 1.1 |
| Naphthalene | ND | | 0.11 |
| 4-Chloroaniline | ND | | 1.1 |
| Hexachlorobutadiene | ND | | 1.1 |
| 4-Chloro-3-methylphenol | ND | | 1.1 |
| 2-Methylnaphthalene | ND | | 0.11 |
| 1-Methylnaphthalene | ND | | 0.11 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 2 of 3

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|------|
| Hexachlorocyclopentadiene | ND | • | 1.1 |
| 2,4,6-Trichlorophenol | ND | | 1.1 |
| 2,3-Dichloroaniline | ND | | 1.1 |
| 2,4,5-Trichlorophenol | ND | | 1.1 |
| 2-Chloronaphthalene | ND | | 1.1 |
| 2-Nitroaniline | ND | | 1.1 |
| 1,4-Dinitrobenzene | ND | | 1.1 |
| Dimethylphthalate | ND | | 1.1 |
| 1,3-Dinitrobenzene | ND | | 1.1 |
| 2,6-Dinitrotoluene | ND | | 1.1 |
| 1,2-Dinitrobenzene | ND | | 1.1 |
| Acenaphthylene | ND | | 0.11 |
| 3-Nitroaniline | ND | | 1.1 |
| 2,4-Dinitrophenol | ND | | 5.4 |
| Acenaphthene | ND | | 0.11 |
| 4-Nitrophenol | ND | | 1.1 |
| 2,4-Dinitrotoluene | ND | | 1.1 |
| Dibenzofuran | ND | | 1.1 |
| 2,3,4,6-Tetrachlorophenol | ND | | 1.1 |
| 2,3,5,6-Tetrachlorophenol | ND | | 1.1 |
| Diethylphthalate | 1.9 | | 1.1 |
| 4-Chlorophenyl-phenylether | ND | | 1.1 |
| 4-Nitroaniline | ND | | 1.1 |
| Fluorene | ND | | 0.11 |
| 4,6-Dinitro-2-methylphenol | ND | | 5.4 |
| N-Nitrosodiphenylamine | ND | | 1.1 |
| 1,2-Diphenylhydrazine | ND | | 1.1 |
| 4-Bromophenyl-phenylether | ND | | 1.1 |
| Hexachlorobenzene | ND | | 1.1 |
| Pentachlorophenol | ND | | 5.4 |
| Phenanthrene | ND | | 0.11 |
| Anthracene | ND | | 0.11 |
| Carbazole | ND | | 1.1 |
| Di-n-butylphthalate | ND | | 1.1 |
| Fluoranthene | ND | | 0.11 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM

page 3 of 3

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|-------|
| Benzidine | ND | | 11 |
| Pyrene | ND | | 0.11 |
| Butylbenzylphthalate | ND | | 1.1 |
| bis-2-Ethylhexyladipate | ND | | 1.1 |
| 3,3'-Dichlorobenzidine | ND | | 1.1 |
| Benzo[a]anthracene | ND | | 0.011 |
| Chrysene | ND | | 0.011 |
| bis(2-Ethylhexyl)phthalate | ND | | 1.1 |
| Di-n-octylphthalate | ND | | 1.1 |
| Benzo[b]fluoranthene | ND | | 0.011 |
| Benzo[k]fluoranthene | ND | | 0.011 |
| Benzo[a]pyrene | ND | | 0.011 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.011 |
| Dibenz[a,h]anthracene | ND | | 0.011 |
| Benzo[g,h,i]perylene | ND | | 0.011 |

| Surrogate : | Percent Recovery | Control Limits |
|----------------------|---------------------|-------------------|
| O Flygraphanal | 40 | 4000 |
| 2-Fluorophenol | 48 | 10 - 90 |
| Phenol-d6 | 37 | 35 - 100 |
| Nitrobenzene-d5 | 67 | 30 - 100 |
| 2-Fluorobiphenyl | 68 | 39 - 100 |
| 2,4,6-Tribromophenol | 82 | 50 - 105 |
| Terphenyl-d14 | 107 | 49 - 115 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 1 of 3

Date Extracted: 3-13-08
Date Analyzed: 3-17&18-08

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0313W1

| Compound: | Results | Flags | PQL |
|---------------------------------|---------|-------|------|
| N-Nitrosodimethylamine | ND | 9- | 1.0 |
| Pyridine | ND | | 1.0 |
| Phenol | ND | | 1.0 |
| Aniline | ND | | 1.0 |
| bis(2-Chloroethyl)ether | ND | | 1.0 |
| 2-Chlorophenol | ND | | 1.0 |
| 1,3-Dichlorobenzene | ND | | 1.0 |
| 1,4-Dichlorobenzene | ND | | 1.0 |
| Benzyl alcohol | ND | | 1.0 |
| 1,2-Dichlorobenzene | ND | | 1.0 |
| 2-Methylphenol (o-Cresol) | ND | | 1.0 |
| bis(2-Chloroisopropyl)ether | ND | | 1.0 |
| (3+4)-Methylphenol (m,p-Cresol) | ND | | 1.0 |
| N-Nitroso-di-n-propylamine | ND | | 1.0 |
| Hexachloroethane | ND | | 1.0 |
| Nitrobenzene | ND | | 1.0 |
| Isophorone | ND | | 1.0 |
| 2-Nitrophenol | ND | | 1.0 |
| 2,4-Dimethylphenol | ND | | 1.0 |
| bis(2-Chloroethoxy)methane | ND | | 1.0 |
| 2,4-Dichlorophenol | ND | | 1.0 |
| 1,2,4-Trichlorobenzene | ND | | 1.0 |
| Naphthalene | ND | | 0.10 |
| 4-Chloroaniline | ND | | 1.0 |
| Hexachlorobutadiene | ND | | 1.0 |
| 4-Chloro-3-methylphenol | ND | | 1.0 |
| 2-Methylnaphthalene | ND | | 0.10 |
| 1-Methylnaphthalene | ND | | 0.10 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 2 of 3

Lab ID: MB0313W1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|------|
| Hexachlorocyclopentadiene | ND | | 1.0 |
| 2,4,6-Trichlorophenol | ND | | 1.0 |
| 2,3-Dichloroaniline | ND | | 1.0 |
| 2,4,5-Trichlorophenol | ND | | 1.0 |
| 2-Chloronaphthalene | ND | | 1.0 |
| 2-Nitroaniline | ND | | 1.0 |
| 1,4-Dinitrobenzene | ND | | 1.0 |
| Dimethylphthalate | ND | | 1.0 |
| 1,3-Dinitrobenzene | ND | | 1.0 |
| 2,6-Dinitrotoluene | ND | | 1.0 |
| 1,2-Dinitrobenzene | ND | | 1.0 |
| Acenaphthylene | ND | | 0.10 |
| 3-Nitroaniline | ND | | 1.0 |
| 2,4-Dinitrophenol | ND | | 5.0 |
| Acenaphthene | ND | | 0.10 |
| 4-Nitrophenol | ND | | 1.0 |
| 2,4-Dinitrotoluene | ND | | 1.0 |
| Dibenzofuran | ND | | 1.0 |
| 2,3,4,6-Tetrachlorophenol | ND | | 1.0 |
| 2,3,5,6-Tetrachlorophenol | ND | | 1.0 |
| Diethylphthalate | ND | | 1.0 |
| 4-Chlorophenyl-phenylether | ND | | 1.0 |
| 4-Nitroaniline | ND | | 1.0 |
| Fluorene | ND | | 0.10 |
| 4,6-Dinitro-2-methylphenol | ND | | 5.0 |
| N-Nitrosodiphenylamine | ND | | 1.0 |
| 1,2-Diphenylhydrazine | ND | | 1.0 |
| 4-Bromophenyl-phenylether | ND | | 1.0 |
| Hexachlorobenzene | ND | | 1.0 |
| Pentachlorophenol | ND | | 5.0 |
| Phenanthrene | ND | | 0.10 |
| Anthracene | ND | | 0.10 |
| Carbazole | ND | | 1.0 |
| Di-n-butylphthalate | ND | | 1.0 |
| Fluoranthene | ND | | 0.10 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM METHOD BLANK QUALITY CONTROL

page 3 of 3

Lab ID: MB0313W1

| Compound: | Results | Flags | PQL |
|----------------------------|---------|-------|-------|
| Benzidine | ND | | 10 |
| Pyrene | ND | | 0.10 |
| Butylbenzylphthalate | ND | | 1.0 |
| bis-2-Ethylhexyladipate | ND | | 1.0 |
| 3,3'-Dichlorobenzidine | ND | | 1.0 |
| Benzo[a]anthracene | ND | | 0.010 |
| Chrysene | ND | | 0.010 |
| bis(2-Ethylhexyl)phthalate | ND | | 1.0 |
| Di-n-octylphthalate | ND | | 1.0 |
| Benzo[b]fluoranthene | ND | | 0.010 |
| Benzo[k]fluoranthene | ND | | 0.010 |
| Benzo[a]pyrene | ND | | 0.010 |
| Indeno[1,2,3-cd]pyrene | ND | | 0.010 |
| Dibenz[a,h]anthracene | ND | | 0.010 |
| Benzo[g,h,i]perylene | ND | | 0.010 |

| Surrogate : | Percent | Control |
|----------------------|----------|----------|
| - | Recovery | Limits |
| 2-Fluorophenol | 47 | 10 - 90 |
| Phenol-d6 | 36 | 35 - 100 |
| Nitrobenzene-d5 | 66 | 30 - 100 |
| 2-Fluorobiphenyl | 65 | 39 - 100 |
| 2,4,6-Tribromophenol | 76 | 50 - 105 |
| Terphenyl-d14 | 97 | 49 - 115 |

Project: 8915 162890

SEMIVOLATILES by EPA 8270D/SIM SB/SBD QUALITY CONTROL

Date Extracted: 3-13-08 Date Analyzed: 3-17-08

Matrix: Water Units: ug/L (ppb)

Lab ID: SB0313W1

| | Spike | | Percent | | Percent I | Recovery | |
|----------------------------|--------|------|----------|------|-----------|----------|-------|
| Compound: | Amount | SB | Recovery | SBD | Recovery | Limits | Flags |
| Phenol | 40.0 | 19.9 | 50 | 20.9 | 52 | 21-75 | |
| | | | | | _ | | |
| 2-Chlorophenol | 40.0 | 35.9 | 90 | 37.1 | 93 | 43-96 | |
| 1,4-Dichlorobenzene | 20.0 | 14.3 | 72 | 15.7 | 78 | 38-80 | |
| N-Nitroso-di-n-propylamine | 20.0 | 17.4 | 87 | 18.2 | 91 | 36-99 | |
| 1,2,4-Trichlorobenzene | 20.0 | 12.8 | 64 | 14.0 | 70 | 39-85 | |
| 4-Chloro-3-methylphenol | 40.0 | 37.5 | 94 | 37.4 | 94 | 50-105 | |
| Acenaphthene | 20.0 | 16.0 | 80 | 16.3 | 82 | 46-90 | |
| 2,4-Dinitrotoluene | 20.0 | 19.3 | 97 | 18.9 | 95 | 50-122 | |
| 4-Nitrophenol | 40.0 | 25.8 | 64 | 25.7 | 64 | 30-116 | |
| Pentachlorophenol | 40.0 | 31.6 | 79 | 31.2 | 78 | 40-112 | |
| Pyrene | 20.0 | 20.1 | 100 | 19.6 | 98 | 51-105 | |

| | RPD | RPD Limits | Flags |
|----------------------------|-----|---------------|-------|
| Phenol | 5 | 31 | |
| 2-Chlorophenol | 3 | 29 | |
| 1,4-Dichlorobenzene | 9 | 33 | |
| N-Nitroso-di-n-propylamine | 4 | 28 | |
| 1,2,4-Trichlorobenzene | 8 | 32 | |
| 4-Chloro-3-methylphenol | 0 | 23 | |
| Acenaphthene | 2 | 24 | |
| 2,4-Dinitrotoluene | 2 | 30 | |
| 4-Nitrophenol | 1 | 30 | |
| Pentachlorophenol | 1 | 30 | |
| Pyrene | 2 | 18 | |

Project: 8915 162890

PCBs by EPA 8082

Matrix: Soil

Units: mg/Kg (ppm)

| | | | | Date | Date | |
|----------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | B-4_3 | | | | | |
| Laboratory ID: | 03-076-01 | | | | | |
| Aroclor 1016 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1221 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1232 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1242 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1248 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1254 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1260 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1262 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1268 | ND | 0.056 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| DCB | 94 | 39-118 | | | | |
| Client ID: | B-5_3 | | | | | |
| Laboratory ID: | 03-076-04 | | | | | |
| Aroclor 1016 | ND | 0.054 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1221 | ND | 0.054 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1232 | ND | 0.054 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1242 | ND | 0.054 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1248 | ND | 0.054 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1254 | ND | 0.054 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1260 | ND | 0.054 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1262 | ND | 0.054 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Aroclor 1268 | ND | 0.054 | EPA 8082 | 3-12-08 | 3-12-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |

Surrogate: Percent Recovery Control Limi
DCB 91 39-118

Project: 8915 162890

PCBs by EPA 8082 QUALITY CONTROL

Matrix: Soil

Units: mg/Kg (ppm)

| | | | Date | Date | |
|----------|---|--|--|--|--|
| Result | PQL | Method | Prepared | Analyzed | Flags |
| | | | | | |
| MB0312S1 | | | | | |
| ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| ND | 0.050 | EPA 8082 | 3-12-08 | 3-12-08 | |
| | MB0312S1 ND ND ND ND ND ND ND ND ND N | MB0312S1 ND 0.050 ND 0.050 | MB0312S1 ND 0.050 EPA 8082 ND 0.050 EPA 8082 | MB0312S1 ND 0.050 EPA 8082 3-12-08 ND 0.050 EPA 8082 3-12-08 | MB0312S1 ND 0.050 EPA 8082 3-12-08 3-12-08 ND 0.050 EPA 8082 3-12-08 3-12-08 |

Surrogate: Percent Recovery Control Limits DCB 101 39-118

| Analyte | Re | sult | Spike | Level | Source Result | | rcent | Recovery Limits | RPD | RPD Limit | Flags |
|----------------|-------|-------|-------|-------|------------------|----|-------|--------------------|-----|--------------|-------|
| MATRIX SPIKES | | | | | | | | | | | |
| Laboratory ID: | 03-0 | 62-05 | | | | | | | | | |
| | MS | MSD | MS | MSD | | MS | MSD | | | | |
| Aroclor 1260 | 0.444 | 0.456 | 0.500 | 0.500 | ND | 89 | 91 | 35-120 | 3 | 17 | |
| Surrogate: | | | | | | | | | | | |
| DCB | | | | | | 93 | 92 | 39-118 | | | |

Project: 8915 162890

ORGANOCHLORINE PESTICIDES by EPA 8081A

Matrix: Soil

Units: ug/Kg (ppb)

| | | | Date | Date | |
|-----------|---|---|---------------------|--|--|
| Result | PQL | Method | Prepared | Analyzed | Flags |
| B-4_3 | | | | | |
| 03-076-01 | | | | | |
| ND | 5.6 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 5.6 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 5.6 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 5.6 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 5.6 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 5.6 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 5.6 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 5.6 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| ND | 56 | EPA 8081 | 3-12-08 | 3-12-08 | |
| | B-4_3 03-076-01 ND | B-4_3 03-076-01 ND 5.6 ND 11 | B-4_3 03-076-01 ND | Result PQL Method Prepared B-4_3 03-076-01 5.6 EPA 8081 3-12-08 ND 11 EPA 8081 3-12-08 N | Result PQL Method Prepared Analyzed B-4_3 03-076-01 3-12-08 3-12-08 3-12-08 ND 5.6 EPA 8081 3-12-08 3-12-08 ND 11 EPA 8081 3-12-08 3-12-08 |

Surrogate: Percent Recovery Control Limits TCMX 74 36-108 DCB 63 30-115

Project: 8915 162890

ORGANOCHLORINE PESTICIDES by EPA 8081A

Matrix: Soil

Units: ug/Kg (ppb)

| 39, 19 (PP2) | | | | Date | Date | |
|--------------------|-----------|-----|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | B-5_3 | | | | | |
| Laboratory ID: | 03-076-04 | | | | | |
| alpha-BHC | ND | 5.4 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-BHC | ND | 5.4 | EPA 8081 | 3-12-08 | 3-12-08 | |
| beta-BHC | ND | 5.4 | EPA 8081 | 3-12-08 | 3-12-08 | |
| delta-BHC | ND | 5.4 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor | ND | 5.4 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Aldrin | ND | 5.4 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor Epoxide | ND | 5.4 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-Chlordane | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| alpha-Chlordane | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDE | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan I | ND | 5.4 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Dieldrin | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDD | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan II | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDT | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Aldehyde | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Methoxychlor | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endsulfan Sulfate | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Ketone | ND | 11 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Toxaphene | ND | 54 | EPA 8081 | 3-12-08 | 3-12-08 | |

Surrogate: Percent Recovery Control Limits TCMX 76 36-108 DCB 76 30-115

Project: 8915 162890

ORGANOCHLORINE PESTICIDES by EPA 8081A QUALITY CONTROL

Matrix: Soil

Units: ug/Kg (ppb)

| Analyte | Result | PQL | Method | Date Propaged | Date Analyzed | Flags |
|--------------------|------------------|----------------|----------|------------------|------------------|-------|
| METHOD BLANK | Result | FQL | Metriou | Prepared | Allalyzeu | riays |
| Laboratory ID: | MB0312S1 | | | | | |
| alpha-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| beta-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| delta-BHC | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Aldrin | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor Epoxide | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-Chlordane | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| alpha-Chlordane | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDE | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan I | ND | 5.0 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Dieldrin | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDD | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan II | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDT | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Aldehyde | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Methoxychlor | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endsulfan Sulfate | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Ketone | ND | 10 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Toxaphene | ND | 50 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| TCMX | 82 | 36-108 | | | | |

Surrogate: Percent Recovery Control Limit
TCMX 82 36-108
DCB 77 30-115

| Analysia | Do | sult | Cnilco | Lovel | Source | | rcent | Recovery Limits | RPD | RPD | Flore |
|----------------|------|-------|--------|-------|--------|-----|-------|--------------------|-----|-------|-------|
| Analyte | Res | Suit | Бріке | Level | Result | Rec | overy | Limits | RPU | Limit | Flags |
| MATRIX SPIKES | | | | | | | | | | | |
| Laboratory ID: | 03-0 | 62-05 | | | | | | | | | |
| | MS | MSD | MS | MSD | | MS | MSD | | | | |
| gamma-BHC | 41.8 | 39.3 | 50.0 | 50.0 | ND | 84 | 79 | 39-106 | 6 | 12 | |
| Heptachlor | 39.3 | 37.3 | 50.0 | 50.0 | ND | 79 | 75 | 33-107 | 5 | 13 | |
| Aldrin | 39.2 | 37.4 | 50.0 | 50.0 | ND | 78 | 75 | 36-101 | 5 | 12 | |
| Dieldrin | 98.1 | 93.0 | 125 | 125 | ND | 78 | 74 | 33-115 | 5 | 11 | |
| Endrin | 99.4 | 94.8 | 125 | 125 | ND | 80 | 76 | 35-108 | 5 | 11 | |
| 4,4'-DDT | 98.4 | 94.3 | 125 | 125 | ND | 79 | 75 | 24-122 | 4 | 17 | |
| Surrogate: | | | | | | | | | | | |
| TCMX | | | | | | 74 | 71 | 36-108 | | | |
| DCB | | | | | | 72 | 70 | 30-115 | | | |

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Project: 8915 162890

ORGANOCHLORINE PESTICIDES by EPA 8081A

Matrix: Water Units: ug/L (ppb)

| | | | | Date | Date | |
|--------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | EB-1 | | | | | |
| Laboratory ID: | 03-076-13 | | | | | |
| alpha-BHC | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-BHC | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| beta-BHC | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| delta-BHC | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Aldrin | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor Epoxide | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-Chlordane | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| alpha-Chlordane | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDE | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan I | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Dieldrin | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDD | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan II | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDT | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Aldehyde | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Methoxychlor | ND | 0.010 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endsulfan Sulfate | ND | 0.0052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Ketone | ND | 0.021 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Toxaphene | ND | 0.052 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Surrogate: | Percent Recovery | Control Limits | · | | | · |

Surrogate: Percent Recovery Control Limit
TCMX 63 30-105
DCB 84 30-121

Project: 8915 162890

ORGANOCHLORINE PESTICIDES by EPA 8081A QUALITY CONTROL

Matrix: Water
Units: ug/L (ppb)

| | | | | Date | Date | |
|--------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| | | | | | | |
| Laboratory ID: | MB0312W1 | | | | | |
| alpha-BHC | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-BHC | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| beta-BHC | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| delta-BHC | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Aldrin | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Heptachlor Epoxide | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| gamma-Chlordane | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| alpha-Chlordane | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDE | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan I | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Dieldrin | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDD | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endosulfan II | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| 4,4'-DDT | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Aldehyde | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Methoxychlor | ND | 0.010 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endsulfan Sulfate | ND | 0.0050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Endrin Ketone | ND | 0.020 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Toxaphene | ND | 0.050 | EPA 8081 | 3-12-08 | 3-12-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| TCMX | 60 | 30-105 | | | | |
| DCB | 88 | 30-121 | | | | |

| | | | | | Source | Pe | rcent | Recovery | | RPD | |
|----------------|--------|--------|--------|--------|--------|-----|-------|----------|-----|-------|-------|
| Analyte | Res | sult | Spike | Level | Result | Rec | overy | Limits | RPD | Limit | Flags |
| SPIKE BLANKS | | | | | | | | | | | |
| Laboratory ID: | SB03 | 12W1 | | | | | | | | | |
| | SB | SBD | SB | SBD | | SB | SBD | | | | |
| gamma-BHC | 0.0374 | 0.0393 | 0.0500 | 0.0500 | N/A | 75 | 79 | 50-104 | 5 | 20 | |
| Heptachlor | 0.0333 | 0.0347 | 0.0500 | 0.0500 | N/A | 67 | 69 | 41-102 | 4 | 22 | |
| Aldrin | 0.0331 | 0.0337 | 0.0500 | 0.0500 | N/A | 66 | 67 | 29-101 | 2 | 25 | |
| Dieldrin | 0.102 | 0.104 | 0.125 | 0.125 | N/A | 81 | 83 | 55-110 | 2 | 24 | |
| Endrin | 0.109 | 0.111 | 0.125 | 0.125 | N/A | 87 | 89 | 54-120 | 2 | 22 | |
| 4,4'-DDT | 0.111 | 0.109 | 0.125 | 0.125 | N/A | 88 | 87 | 56-119 | 2 | 26 | |
| Surrogate: | | | | | | | | | | | |
| TCMX | | | | | | 58 | 58 | 30-105 | | | |
| DCB | | | | | | 82 | 82 | 30-121 | | | |

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Project: 8915 162890

CHLORINATED ACID HERBICIDES by EPA 8151A

Matrix: Soil

Dinoseb

Units: ug/Kg (ppb)

| | | | | Date | Date | |
|-------------------|------------------|----------------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | B-4_3 | | | | | |
| Laboratory ID: | 03-076-01 | | | | | |
| Dalapon | ND | 260 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 5300 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 5300 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 1.1 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-T | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-DB | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dinoseb | ND | 53 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Surrogate: | Percent Recovery | Control Limits | | | | |
| DCAA | 74 | 37-114 | | | | |
| Client ID. | D.E. 2 | | | | | |
| Client ID: | B-5_3 | | | | | |
| Laboratory ID: | 03-076-04 | | | | | |
| Dalapon | ND | 250 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 51 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 5000 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 5000 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 51 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 51 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 1.0 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 51 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-T | ND | 51 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-DB | ND | 51 | EPA 8151 | 3-18-08 | 3-18-08 | |

EPA 8151

3-18-08

3-18-08

Surrogate: Percent Recovery Control Limits DCAA 79 37-114

51

Project: 8915 162890

CHLORINATED ACID HERBICIDES by EPA 8151A QUALITY CONTROL

Matrix: Soil

Units: ug/Kg (ppb)

| | | | | Date | Date | |
|-------------------|----------|------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| METHOD BLANK | | | | | | |
| Laboratory ID: | MB0318S1 | | | | | |
| Dalapon | ND | 230 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dicamba | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPP | ND | 4700 | EPA 8151 | 3-18-08 | 3-18-08 | |
| MCPA | ND | 4700 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dichlorprop | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-D | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Pentachlorophenol | ND | 0.95 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-TP (Silvex) | ND | 48 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4,5-T | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| 2,4-DB | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| Dinoseb | ND | 47 | EPA 8151 | 3-18-08 | 3-18-08 | |
| | | | | | | |

Surrogate: Percent Recovery Control Limits DCAA 70 37-114

| Analyte | Re | sult | Spike | Level | Source Result | | rcent | Recovery Limits | RPD | RPD Limit | Flags |
|----------------|------|-------|-------|-------|------------------|----|-------|--------------------|-----|--------------|-------|
| MATRIX SPIKES | | | | | | | | | | | |
| Laboratory ID: | 03-0 | 62-05 | | | | | | | | | |
| | MS | MSD | MS | MSD | | MS | MSD | | | | |
| Dicamba | 77.7 | 76.9 | 100 | 100 | ND | 78 | 77 | 62-99 | 1 | 10 | |
| 2,4-D | 71.5 | 69.1 | 100 | 100 | ND | 72 | 69 | 23-111 | 3 | 24 | |
| 2,4,5-T | 79.5 | 75.5 | 100 | 100 | ND | 79 | 76 | 38-120 | 5 | 12 | |
| 2,4-DB | 90.3 | 83.9 | 100 | 100 | ND | 90 | 84 | 44-135 | 7 | 16 | |
| Surrogate: | • | • | • | • | | | • | | • | | |
| DCAA | | | | | | 80 | 79 | 37-114 | | | |

Project: 8915 162890

CHLORINATED ACID HERBICIDES by EPA 8151A

Matrix: Water
Units: ug/L (ppb)

| | | | | Date | Date | |
|-------------------|-----------|--------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| Client ID: | EB-1 | | | | | |
| Laboratory ID: | 03-076-13 | | | | | |
| Dalapon | ND | 0.23 | EPA 8151 | 3-13-08 | 3-14-08 | |
| Dicamba | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| MCPP | ND | 4.8 | EPA 8151 | 3-13-08 | 3-14-08 | |
| MCPA | ND | 4.8 | EPA 8151 | 3-13-08 | 3-14-08 | |
| Dichlorprop | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| 2,4-D | ND | 0.48 | EPA 8151 | 3-13-08 | 3-14-08 | |
| Pentachlorophenol | ND | 0.0097 | EPA 8151 | 3-13-08 | 3-14-08 | |
| 2,4,5-TP (Silvex) | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| 2,4,5-T | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| 2,4-DB | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| Dinoseb | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |

Surrogate: Percent Recovery Control Limits DCAA 102 44-116

Project: 8915 162890

CHLORINATED ACID HERBICIDES by EPA 8151A QUALITY CONTROL

Matrix: Water
Units: ug/L (ppb)

| | | | | Date | Date | |
|-------------------|----------|--------|----------|----------|----------|-------|
| Analyte | Result | PQL | Method | Prepared | Analyzed | Flags |
| METHOD BLANK | | | | | | |
| Laboratory ID: | MB0313W1 | | | | | |
| Dalapon | ND | 0.23 | EPA 8151 | 3-13-08 | 3-14-08 | |
| Dicamba | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| MCPP | ND | 4.7 | EPA 8151 | 3-13-08 | 3-14-08 | |
| MCPA | ND | 4.7 | EPA 8151 | 3-13-08 | 3-14-08 | |
| Dichlorprop | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| 2,4-D | ND | 0.47 | EPA 8151 | 3-13-08 | 3-14-08 | |
| Pentachlorophenol | ND | 0.0095 | EPA 8151 | 3-13-08 | 3-14-08 | |
| 2,4,5-TP (Silvex) | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| 2,4,5-T | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| 2,4-DB | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| Dinoseb | ND | 0.024 | EPA 8151 | 3-13-08 | 3-14-08 | |
| <u> </u> | | | · | | | |

Surrogate: Percent Recovery Control Limits DCAA 97 44-116

| | | | | | Source | Pe | rcent | Recovery | | RPD | |
|----------------|-------|-------|-------|-------|--------|-----|-------|----------|-----|-------|-------|
| Analyte | Re | sult | Spike | Level | Result | Rec | overy | Limits | RPD | Limit | Flags |
| SPIKE BLANKS | | | | | | | | | | | |
| Laboratory ID: | SB03 | 13W1 | | | | | | | | | |
| | SB | SBD | SB | SBD | | SB | SBD | | | | |
| Dicamba | 0.830 | 0.854 | 1.00 | 1.00 | N/A | 83 | 85 | 29-121 | 3 | 20 | _ |
| 2,4-D | 0.842 | 0.858 | 1.00 | 1.00 | N/A | 84 | 86 | 47-110 | 2 | 20 | |
| 2,4,5-T | 0.912 | 0.930 | 1.00 | 1.00 | N/A | 91 | 93 | 61-119 | 2 | 20 | |
| 2,4-DB | 0.992 | 1.01 | 1.00 | 1.00 | N/A | 99 | 101 | 70-144 | 2 | 20 | |
| Surrogate: | • | | • | | • | | • | | | | |
| DCAA | | | | | | 97 | 102 | 44-116 | | | |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A

Date Extracted: 3-13&14-08 Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-076-03 **Client ID: B-4_8**

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 6010B | ND | 11 |
| Barium | 6010B | 33 | 2.7 |
| Cadmium | 6010B | ND | 0.55 |
| Chromium | 6010B | 18 | 0.55 |
| Lead | 6010B | ND | 5.5 |
| Mercury | 7471A | ND | 0.27 |
| Selenium | 6010B | ND | 11 |
| Silver | 6010B | ND | 0.55 |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A

Date Extracted: 3-13&14-08 Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-076-09 **Client ID: B-5_11**

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 6010B | ND | 11 |
| Barium | 6010B | 34 | 2.9 |
| Cadmium | 6010B | ND | 0.57 |
| Chromium | 6010B | 26 | 0.57 |
| Lead | 6010B | ND | 5.7 |
| Mercury | 7471A | ND | 0.29 |
| Selenium | 6010B | ND | 11 |
| Silver | 6010B | ND | 0.57 |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A

Date Extracted: 3-13&14-08 Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-076-10 Client ID: B-5_12

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 6010B | ND | 11 |
| Barium | 6010B | 39 | 2.8 |
| Cadmium | 6010B | ND | 0.57 |
| Chromium | 6010B | 30 | 0.57 |
| Lead | 6010B | ND | 5.7 |
| Mercury | 7471A | ND | 0.28 |
| Selenium | 6010B | ND | 11 |
| Silver | 6010B | ND | 0.57 |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A METHOD BLANK QUALITY CONTROL

Date Extracted: 3-13&14-08
Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0313S2&MB0314S2

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 6010B | ND | 10 |
| Barium | 6010B | ND | 2.5 |
| Cadmium | 6010B | ND | 0.50 |
| Chromium | 6010B | ND | 0.50 |
| Lead | 6010B | ND | 5.0 |
| Mercury | 7471A | ND | 0.25 |
| Selenium | 6010B | ND | 10 |
| Silver | 6010B | ND | 0.50 |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A DUPLICATE QUALITY CONTROL

Date Extracted: 3-13&14-08 Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-087-19

| Analyte | Sample Result | Duplicate Result | RPD | PQL | Flags |
|----------|------------------|---------------------|-----|------|-------|
| Arsenic | 13.7 | 13.0 | 5 | 10 | |
| Barium | 87.8 | 84.6 | 4 | 2.5 | |
| Cadmium | ND | ND | NA | 0.50 | |
| Chromium | 35.0 | 33.0 | 6 | 0.50 | |
| Lead | 13.5 | 15.1 | 12 | 5.0 | |
| Mercury | ND | ND | NA | 0.25 | |
| Selenium | ND | ND | NA | 10 | |
| Silver | ND | ND | NA | 0.50 | |

Project: 8915 162890

TOTAL METALS EPA 6010B/7471A MS/MSD QUALITY CONTROL

Date Extracted: 3-13&14-08 Date Analyzed: 3-13&17-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-087-19

| Analyte | Spike Level | MS | Percent Recovery | MSD | Percent Recovery | RPD | Flags |
|----------|----------------|-------|---------------------|-------|---------------------|-----|-------|
| Arsenic | 100 | 103 | 90 | 105 | 91 | 2 | |
| Barium | 100 | 185 | 97 | 193 | 105 | 5 | |
| Cadmium | 50 | 47.6 | 95 | 48.2 | 96 | 1 | |
| Chromium | 100 | 129 | 94 | 128 | 93 | 0 | |
| Lead | 250 | 244 | 92 | 267 | 102 | 9 | |
| Mercury | 0.50 | 0.494 | 99 | 0.504 | 101 | 2 | |
| Selenium | 100 | 87.5 | 87 | 88.3 | 88 | 1 | |
| Silver | 25 | 21.9 | 88 | 22.1 | 88 | 1 | |

Project: 8915 162890

TOTAL METALS EPA 200.8/7470A

Date Extracted: 3-17&19-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-076-13 **Client ID: EB-1**

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 200.8 | ND | 3.3 |
| Barium | 200.8 | ND | 28 |
| Cadmium | 200.8 | ND | 4.4 |
| Chromium | 200.8 | ND | 11 |
| Lead | 200.8 | ND | 1.1 |
| Mercury | 7470A | ND | 0.50 |
| Selenium | 200.8 | ND | 5.6 |
| Silver | 200.8 | ND | 11 |

Project: 8915 162890

TOTAL METALS EPA 200.8/7470A METHOD BLANK QUALITY CONTROL

Date Extracted: 3-17&19-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: MB0317W1&MB0319W1

| Analyte | Method | Result | PQL |
|----------|--------|--------|------|
| Arsenic | 200.8 | ND | 3.3 |
| Barium | 200.8 | ND | 28 |
| Cadmium | 200.8 | ND | 4.4 |
| Chromium | 200.8 | ND | 11 |
| Lead | 200.8 | ND | 1.1 |
| Mercury | 7470A | ND | 0.50 |
| Selenium | 200.8 | ND | 5.6 |
| Silver | 200.8 | ND | 11 |

Project: 8915 162890

TOTAL METALS EPA 200.8/7470A DUPLICATE QUALITY CONTROL

Date Extracted: 3-17&19-08 Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-076-13

| Analyte | Sample Result | Duplicate Result | RPD | PQL | Flags |
|----------|------------------|---------------------|-----|------|-------|
| Arsenic | ND | ND | NA | 3.3 | |
| Barium | ND | ND | NA | 28 | |
| Cadmium | ND | ND | NA | 4.4 | |
| Chromium | ND | ND | NA | 11 | |
| Lead | ND | ND | NA | 1.1 | |
| Mercury | ND | ND | NA | 0.50 | |
| Selenium | ND | ND | NA | 5.6 | |
| Silver | ND | ND | NA | 11 | |

Project: 8915 162890

TOTAL METALS EPA 200.8/7470A MS/MSD QUALITY CONTROL

Date Extracted: 3-17&19-08
Date Analyzed: 3-18&19-08

Matrix: Water Units: ug/L (ppb)

Lab ID: 03-076-13

| Analyte | Spike Level | MS | Percent Recovery | MSD | Percent Recovery | RPD | Flags |
|----------|----------------|------|---------------------|-----|---------------------|-----|-------|
| Arsenic | 110 | 113 | 102 | 113 | 103 | 0 | |
| Barium | 110 | 116 | 106 | 115 | 104 | 1 | |
| Cadmium | 110 | 114 | 103 | 119 | 108 | 4 | |
| Chromium | 110 | 112 | 102 | 113 | 103 | 1 | |
| Lead | 110 | 112 | 102 | 112 | 102 | 0 | |
| Mercury | 12.5 | 11.8 | 94 | 12 | 96 | 2 | |
| Selenium | 110 | 119 | 109 | 116 | 105 | 3 | |
| Silver | 110 | 108 | 98 | 113 | 103 | 5 | |

Project: 8915 162890

SOLUBLE HEXAVALENT CHROMIUM WATER EXTRACTION EPA 7196A

Date Extracted: 3-24-08
Date Analyzed: 3-24-08

Matrix: Soil

Units: mg/kg (ppm)

| Client ID | Lab ID | Result | PQL |
|-----------|-----------|--------|-----|
| B-4_8 | 03-076-03 | ND | 1.1 |
| B-5_11 | 03-076-09 | ND | 1.1 |
| B-5_12 | 03-076-10 | ND | 1.1 |

Project: 8915 162890

SOLUBLE HEXAVALENT CHROMIUM WATER EXTRACTION EPA 7196A METHOD BLANK QUALITY CONTROL

Date Extracted: 3-24-08 Date Analyzed: 3-24-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: MB0324S1

Analyte Method Result PQL
Hexavalent Chromium 7196A **ND** 1.0

Project: 8915 162890

SOLUBLE HEXAVALENT CHROMIUM WATER EXTRACTION EPA 7196A DUPLICATE QUALITY CONTROL

Date Extracted: 3-24-08 Date Analyzed: 3-24-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-05

Sample Duplicate

Analyte Result Result RPD Flags PQL
Hexavalent Chromium ND ND NA 1.0

Project: 8915 162890

SOLUBLE HEXAVALENT CHROMIUM WATER EXTRACTION EPA 7196A MS/MSD QUALITY CONTROL

Date Extracted: 3-24-08 Date Analyzed: 3-24-08

Matrix: Soil

Units: mg/kg (ppm)

Lab ID: 03-062-05

| Analyte | Spike Level | MS | Percent Recovery | MSD | Percent Recovery | RPD | Flags |
|---------------------|----------------|------|---------------------|------|---------------------|-----|-------|
| Hexavalent Chromium | 5.0 | 4.76 | 95 | 4.57 | 91 | 4 | |

Project: 8915 162890

% MOISTURE

Date Analyzed: 3-12-08

| Client ID | Lab ID | % Moisture |
|-----------|-----------|------------|
| B_4_3 | 03-076-01 | 11 |
| B-4_6 | 03-076-02 | 11 |
| B-4_8 | 03-076-03 | 9 |
| B-5_3 | 03-076-04 | 7 |
| B-5_6 | 03-076-05 | 10 |
| B-5_7 | 03-076-06 | 9 |
| B-5_8 | 03-076-07 | 4 |
| B-5_9 | 03-076-08 | 4 |
| B-5_11 | 03-076-09 | 13 |
| B-5_12 | 03-076-10 | 12 |



Data Qualifiers and Abbreviations

- A Due to a high sample concentration, the amount spiked is insufficient for meaningful MS/MSD recovery data.
- B The analyte indicated was also found in the blank sample.
- C The duplicate RPD is outside control limits due to high result variability when analyte concentrations are within five times the quantitation limit.
- E The value reported exceeds the quantitation range and is an estimate.
- F Surrogate recovery data is not available due to the high concentration of coeluting target compounds.
- H The analyte indicated is a common laboratory solvent and may have been introduced during sample preparation, and be impacting the sample result.
- I Compound recovery is outside of the control limits.
- J The value reported was below the practical quantitation limit. The value is an estimate.
- K Sample duplicate RPD is outside control limits due to sample inhomogeneity. The sample was re-extracted and re-analyzed with similar results.
- L The RPD is outside of the control limits.
- M Hydrocarbons in the gasoline range are impacting the diesel range result.
- M1 Hydrocarbons in the gasoline range (toluene-napthalene) are present in the sample.
- N Hydrocarbons in the lube oil range are impacting the diesel range result.
- O Hydrocarbons indicative of heavier fuels are present in the sample and are impacting the gasoline result.
- P The RPD of the detected concentrations between the two columns is greater than 40.
- Q Surrogate recovery is outside of the control limits.
- S Surrogate recovery data is not available due to the necessary dilution of the sample.
- T The sample chromatogram is not similar to a typical ______
- U The analyte was analyzed for, but was not detected above the reported sample quantitation limit.
- U1 The practical quantitation limit is elevated due to interferences present in the sample.
- V Matrix Spike/Matrix Spike Duplicate recoveries are outside control limits due to matrix effects.
- W Matrix Spike/Matrix Spike Duplicate RPD are outside control limits due to matrix effects.
- X Sample extract treated with a mercury cleanup procedure.
- Y Sample extract treated with an acid/silica gel cleanup procedure.
- Z Analyte may have been introduced by an outside source. See case narrative.
- ND Not Detected at PQL
- PQL Practical Quantitation Limit
- RPD Relative Percent Difference

Chain of Custody

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% Moisture ex chang 8 HEM by 1664 TCLP Metals Laboratory Number: 03-076 > <Total RCRA Metals (8) × Herbicides by 8151A Pesticides by 8081A PCBs by 8082 MIS \ G07S8 vd sHA9 × × Semivolatiles by 8270D Halogenated Volatiles by 8260B X 4008 Volatiles by 8260B **XQ-H9TWN** X378/x2-H9TWN имтрн-нсір (TPH analysis 5 working days) ☐ 1 Day 🗌 з Dау # of Cont. V Standard (7 working days) Turnaround Request (in working days) (Check One) (other) 800 8/0 1220 1200 1225 820 0/2/0 Same Day 2 Day 3/1/8 Date Sampled 100 OnSite Environmental Inc. Phone: (425) 883-3881 • Fax: (425) 885-4603 Manster Sta hen byn Inaye Sample Identification Project Number: 8915 3 00 8 Sampled by: Company: Lab ID

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| Signature | Company Date Time | Comments/Special Instructions: |
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Chain of Custody

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